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C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\2323tg.str
```

```
ring nodes :
   1 2 3
                       8
                          9 10
                                 11
                                    22
                                                    26
chain bonds :
                                                30-33
   6-22 12-13
               23-27
                      25-28
                             28-29
                                   29-30
                                          30-31
ring bonds :
   1-2 1-6 2-3 2-7
                      3-4 3-10 4-5 5-6
                                         7-8 8-11 9-11
                                                          9-10
   22-26 23-24 24-25 25-26
exact/norm bonds :
                                                             23-24
   2-7 3-10 6-22 7-8 8-11
                                                      22-26
                              9-11 9-10 12-13 22-23
                                                                    23-27
   24-25 25-26 25-28 28-29 29-30 30-31 30-33
normalized bonds :
   1-2 1-6 . 2-3 3-4 4-5 5-6
isolated ring systems :
   containing 1 : 22 :
G1: [*1], [*2], [*3], [*4]
G2:CH3,Et
Match level :
   1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom
                                                 7:Atom 8:Atom 9:Atom
                                                15:CLASS 16:CLASS 22:Atom
   10:Atom 11:Atom 12:CLASS 13:CLASS 14:CLASS
   23:Atom 24:Atom 25:Atom 26:Atom 27:CLASS 28:CLASS 29:CLASS 30:CLASS
   31:CLASS 33:CLASS
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30 . 31

33

chain nodes :

12 13 14

16

15

27

28

29

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C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\55651.str
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```
ring nodes :
   1 2 3 4
               5 .
chain bonds :
   1-18 2-6
             4-7 7-15
ring bonds :
   1-2 1-5 2-3 3-4 4-5
exact/norm bonds :
   1-2 1-5 1-18 2-3 2-6 3-4 4-5 7-15
exact bonds :
   4-7
G1:[*1],[*2],[*3]
Match level :
   1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:CLASS 8:CLASS 9:CLASS
   10:CLASS 15:CLASS 18:Atom
Generic attributes :
   18:
                      : Unsaturated
   Saturation
Element Count :
   Node 18: Limited
       C, C3-11
       N, NO-4
       0,00-3
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chain nodes :

6 7 8 9 10 15

18

L23 ANSWER 21 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

2002:575073 HCAPLUS

DOCUMENT NUMBER:

137:140512

TITLE:

Preparation of benzoheterocyclyloxazolidinones as

antibacterial agents.

INVENTOR(S):

Johnson, Paul D.; Aristoff, Paul A.; Poel, Toni-Jo;

Thomasco, Lisa M.

PATENT ASSIGNEE(S):

Pharmacia & Upjohn Company, USA

SOURCE:

PCT Int. Appl., 95 pp. CODEN: PIXXD2

DOCUMENT TYPE:

Patent

LANGUAGE:

English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

	PATENT NO.								APPLICATION NO.						DATE			
	2002	0591	15		A1		2002	0801		WO 2	001-	US42	944		. 2	0011	114	
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		CO.	CR.	CU.	CZ.	DE,	DK,	DM,	DZ,	EC,	EE,	ES,	FI,	GB,	GD,	GE,	GH,	
		GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KΕ,	KG,	KP,	KR,	ΚZ,	LC,	LK,	LR,	
		LS.	LT.	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	ΜZ,	NO,	ΝZ,	OM,	PH,	
		PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	ТJ,	TM,	TR,	TT,	ΤZ,	UA,	
		UG,	US,	UZ,	VN,	YU,	ZA,	ZW										
	RW:	GH,	GM,	ΚE,	LS,	MW,	ΜZ,	SĐ,	SL,	SZ,	ΤZ,	UG,	ZW,	ΑT,	BE,	CH,	CY,	
		DE,	DK,	ES,	FI,	FR,	GB,	GR,	ΙE,	IT,	LU,	MC,	NL,	PΤ,	SE,	TR,	BF,	
		ВJ,	CF,	CG,	CI,	CM,	GA,	GN,	GQ,	GW,	ML,	MR,	ΝE,	SN,	TD,	ΤG		
CA	2421	583			AA		2002	0801		CA 2	001-	2421	583.		2	0011	114	
US	2002	1330	21		A1		2002	0919		US 2	001-	9926	60		2	0011	114	
US	6972	286			B2		2005	1206							_			
EP	1337	530			A1		2003	0827		EP 2	001-	9970	07		2	0011	114	
	R:						ES,					LI,	LU,	ΝL,	SE,	MC,	PT,	
		ΙE,	SI,	LT,	LV,	FI,	RO,	MK,	CY,	ΑL,	TR				_			
									JP 2002-559417						20011114			
NZ	NZ 525918				A 20051125				NZ 2001-525918 US 2004-804389						20011114			
US	2004	1766	09		A1		2004	0909		US ₂	004-	8043	89		2	0040	319	
ÜS	2004	1862	93		A1		2004	0923		US 2	004-	8043	80		_ 2	0040	319	
PRIORIT	RIORITY APPLN. INFO.:									US 2	000-	2495	50P		P 2	0001	11/	
										US 2								
										WO 2	001-	US42	944		W 2	0011	114	
OTHER S	OTHER SOURCE(S):				MAR	PAT	137:	1405	12									

(%) (%)

AB Title compds. [I; Y = NHC(:W)R1, OZ, SZ, NHZ; X = O, NR3, S, SO, SO2, S(O)(NR4); W = O, S; R1 = H, alkyl, cycloalkyl, alkoxy, alkylthio, amino;

Updated Search

GI

R2 = H, halo, alkyl; R3 = H, alkyl, aryl, Z, etc.; R4 = H, alkyl; Z = (substituted) (aromatic) heterocyclyl; with provisos], were prepared Thus, N-[[(5S)-3-(1,2,3,4-tetrahydro-6-isoquinolinyl)-2-oxo-5-oxazolidinyl]methyl]acetamide (preparation given) and NaHCO3 in THF were treated with MeO2CCl with vigorous stirring. H2O was added and the mixture was stirred 1 h to give Me (-)-6-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3,4-dihydro-2(1H)-isoquinolinecarboxylate. The latter showed a min. inhibitory concentration of 1 μ g/mL against Streptococcus pneumoniae SPNE9912.

IT 444588-30-1P RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(preparation of benzoheterocyclyloxazolidinones as antibacterial agents)

RN 444588-30-1 HCAPLUS

CN 2H-2-Benzazepine-2-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,3,4,5-tetrahydro-, phenylmethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

```
444587-37-5P 444587-38-6P 444587-39-7P
IT
     444587-40-0P 444587-46-6P 444587-48-8P
     444587-52-4P 444587-55-7P 444587-56-8P
     444587-57-9P 444587-58-0P 444587-60-4P
     444587-61-5P 444587-62-6P 444587-63-7P
     444587-64-8P 444587-65-9P 444587-66-0P
     444587-67-1P 444587-68-2P 444587-69-3P
     444587-70-6P 444587-71-7P 444587-72-8P
     444587-73-9P 444587-74-0P 444587-75-1P
     444587-76-2P 444587-77-3P 444587-78-4P
     444587-79-5P 444587-80-8P 444587-81-9P
     444587-83-1P 444587-84-2P 444587-85-3P
     RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
     (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
     (Uses)
```

(preparation of benzoheterocyclyloxazolidinones as antibacterial agents) 444587-37-5 HCAPLUS

2(1H)-Isoquinolinecarboxylic acid, 6-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3,4-dihydro-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN

CN

RN 444587-38-6 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-formyl-1,2,3,4-tetrahydro-6-isoquinolinyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-39-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-[2-[(acetyloxy)acetyl]-1,2,3,4-tetrahydro-6-isoquinolinyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-40-0 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[1,2,3,4-tetrahydro-2-(hydroxyacetyl)-6-isoquinolinyl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444587-46-6 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-1H-2-benzopyran-6-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-48-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-1H-2-benzothiopyran-6-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-52-4 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-1H-2-benzothiopyran-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-55-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-2-oxido-1H-2-benzothiopyran-7-y1)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-56-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3-formyl-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-57-9 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-(hydroxyacetyl)-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-58-0 HCAPLUS

CN 3H-3-Benzazepine-3-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4,5-tetrahydro-, phenylmethyl ester (9CI) (CA INDEX NAME)

外

RN 444587-60-4 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3-acetyl-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



RN 444587-61-5 HCAPLUS

CN 3H-3-Benzazepine-3-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4,5-tetrahydro-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-62-6 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3-benzoyl-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

RN 444587-63-7 HCAPLUS
CN Acetamide, N-[[(5S)-3-[3-(5-amino-1,3,4-thiadiazol-2-yl)-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-64-8 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-(methylsulfonyl)-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-65-9 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-[5-(methylthio)-1,3,4-thiadiazol-2-yl]-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444587-66-0 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-(5-methyl-1,3,4-thiadiazol-2-yl)-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-67-1 HCAPLUS

CN 3H-3-Benzazepine-3-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4,5-tetrahydro-, phenyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-68-2 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-(phenylacetyl)-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444587-69-3 HCAPLUS
CN Acetamide, N-[[(5S)-3-[3-[5-(formylamino)-1,3,4-thiadiazol-2-yl]-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-70-6 HCAPLUS
CN Acetamide, N-[5-[7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]1,2,4,5-tetrahydro-3H-3-benzazepin-3-yl]-1,3,4-thiadiazol-2-yl]-2-hydroxy(9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-71-7 HCAPLUS
CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-[(4-iodophenyl)acetyl]1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444587-72-8 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-[[3-(trifluoromethyl)phenyl]acetyl]-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-73-9 HCAPLUS

CN Benzoic acid, 4-[(dimethylamino)methyl]-, 2-[7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4,5-tetrahydro-3H-3benzazepin-3-yl]-2-oxoethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-74-0 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-3-[[4-(trifluoromethyl)phenyl]acetyl]-1H-3-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-75-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3-(1,4-dioxopentyl)-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-76-2 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3-(1,5-dioxohexyl)-2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-77-3 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-formyl-2,3,4,5-tetrahydro-1H-2-benzazepin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444587-78-4 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2,3,4,5-tetrahydro-2-(hydroxyacetyl)-1H-2-benzazepin-7-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-79-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-acetyl-2,3,4,5-tetrahydro-1H-2-benzazepin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-80-8 HCAPLUS

CN 2H-2-Benzazepine-2-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,3,4,5-tetrahydro-, methyl ester (9CI) (CA INDEX NAME)

RN 444587-81-9 HCAPLUS

CN 2H-2-Benzazepine-2-carboxamide, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,3,4,5-tetrahydro-N-phenyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 444587-83-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-(1-formyl-2,3,4,5-tetrahydro-1H-1-benzazepin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

A

RN 444587-84-2 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,4,5-tetrahydro-3-benzothiepin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444587-85-3 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,4,5-tetrahydro-3,3-dioxido-3-benzothiepin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 444587-91-1P 444587-92-2P 444588-02-7P

444588-04-9P 444588-08-3P 444588-20-9P

444588-21-0P 444588-27-6P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of benzoheterocyclyloxazolidinones as antibacterial agents)

RN 444587-91-1 HCAPLUS

CN 2(1H)-Isoquinolinecarboxylic acid, 6-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3,4-dihydro-, phenylmethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444587-92-2 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-6-isoquinolinyl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 444588-02-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-2,2-dioxido-1H-2-benzothiopyran-6-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

444588-04-9 HCAPLUS

RN

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-1H-2-benzopyran-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444588-08-3 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-2,2-dioxido-1H-2-benzothiopyran-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444588-20-9 HCAPLUS

CN 1H-1-Benzazepine-1-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,3,4,5-tetrahydro-, phenylmethyl ester (9CI) (CA INDEX

Updated Search

NAME)

Absolute stereochemistry.

RN 444588-21-0 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(2,3,4,5-tetrahydro-1H-1-benzazepin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 444588-27-6 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(2,3,4,5-tetrahydro-1H-3-benzazepin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

C:\Documents and Settings\brobinson1\My Documents\stnweb\Queries\lmlmp.str chain nodes : 12 1 8 9 10 11

```
10-11
                             11-12
   1-3 4-8 6-9 9-10
ring bonds :
   3-4 3-7
            4-5
exact/norm bonds :
                             11-12
                  4 – 8
                      10-11
                                    11-14
   1-3 3-4 3-7
exact bonds :
   4-5 5-6 6-7 6-9
                      9-10
isolated ring systems :
   containing 3 :
G1
G2:CH3,Et
Match level :
   1:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:CLASS 9:CLASS 10:CLASS
   11:CLASS 12:CLASS 14:CLASS
Generic attributes :
   1:
                         : Unsaturated
   Saturation
   Number of Carbon Atoms: 7 or more
   Type of Ring System : Polycyclic
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ring nodes :

chain bonds :

3 4 5 6

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 29 OF 50

ACCESSION NUMBER:

1999:511162 HCAPLUS

DOCUMENT NUMBER:

131:144608

TITLE:

Preparation of New oxazolidinones with azo-containing

tricycles as antimicrobial agents

INVENTOR(S):

Raddatz, Siegfried; Bartel, Stephan; Guarnieri, Walter; Rosentreter, Ulrich; Ruppelt, Martin; Wild,

Hanno; Endermann, Rainer; Kroll, Hein-peter;

Henninger, Kerstin

PATENT ASSIGNEE(S):

Bayer Aktiengesellschaft, Germany

SOURCE:

PCT Int. Appl., 207 pp. CODEN: PIXXD2

DOCUMENT TYPE:

Patent

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PA	TENT :	NO.			KIN	D	DATE				ICAT:				D.	ATE	
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		KE,	KG,	ΚP,	KR,	ΚZ,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	MD,	MG,	MK,	MN,
		MW,	MX,	NO,	ΝZ,	PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	ТJ,	TM,
		TR,	TT,	UA,	UG,	US,	UZ,	VN,	YU,	ZW,	AM,	ΑZ,	BY,	KG,	ΚZ,	MD,	RU,
		ТJ,												•			
	RW:	GH,	GM,	ΚE,	LS,	MW,	SD,	SZ,	UG,	ZW,	ΑT,	BE,	CH,	CY,	DE,	DK,	ES,
											PT,	SE,	BF,	·BJ,	CF,	CG,	CI,
		CM,	GA,	GN,	GW,	ML,	MR,	NE,	SN,	TD,	TG						
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EP	1054																
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		ΙE,															
JP	2002	5028	54		Т2		2002	0129								9990	
PRIORIT	Y APP	LN.	INFO	.:							998-						
										WO 1	999-	EP51	8	,	W 1	9990	127
OTHER S	OURCE	(S):			MAR	PAT	131:	1446	80	•							

The present invention relates to new oxazolidinones with azo-containing tricycles, to methods for producing the same as well as to the use thereof as drugs, in particular as anti-bacterial drugs. Title compds. [I; R = NHAC, NHCOCH2C1, NHCOOME, NH2, NHCOCHC12, NHCOEt, NHCOCC13, NHCSNH2, NHCSNHME, NHCONH2, NHCHO, NHCSME, NHCOOBu-t, NHCSOME, NHCSOEt, OH, 3-NO2C6H4SO2O; A = O, CH2, S; B = CH2, (CH3)2C, S, O, S:O, SO2, CF2; B-A = S, CH2, CH:CH; X = CH, CCH3, N, CBr; Y = N, CCH3, CPh, CCOOEt, CCF3, CNHCOOBu-t, CBr, CCH2OH; Z = N, CCOOEt, CH], enantiomers, and salts are prepared as antibacterial drugs, tested against Staphylococcus aureus, Mycobacterium smegmatis, and Streptococcus pneumoniae. Thus, the title compound II was prepared from (5S)-3-(2-aminobenzthiazol-6-yl)-5-acetylaminomethyl-oxazolidin-2-one and CH2ClCHO via cyclization and tested.

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TT 235789-77-2P 235789-78-3P 235789-79-4P 235789-80-7P 235789-81-8P 235789-82-9P 235789-83-0P 235789-85-2P 235789-86-3P 235789-87-4P 235789-88-5P 235789-89-6P 235789-90-9P 235789-91-0P 235789-92-1P 235789-93-2P 235789-94-3P 235789-95-4P 235789-98-8P 235789-97-6P 235789-98-7P 235789-98-P 235790-01-9P 235790-02-0P 235790-03-1P 235790-04-2P 235790-17-7P 235790-18-8P 235790-19-9P 235790-26-8P 235790-27-9P 235792-17-3P RI: BAC (Biological activity or effector. Expenses the second content of the second cont
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RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of oxazolidines substituted with bicycles as antimicrobial agents)

RN 235789-77-2 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(4H-[1,2,4]triazolo[3,4-c][1,4]benzoxazin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235789-78-3 HCAPLUS

CN Propanamide, N-[[(5S)-3-(1-methyl-4H-[1,2,4]triazolo[3,4-c][1,4]benzoxazin-7-y1)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-79-4 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4,5-dihydro[1,2,4]triazolo[4,3-a]quinolin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-80-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(1-methyl-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235789-81-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-methyl-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-82-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-83-0 HCAPLUS

CN Propanamide, N-[[(5S)-3-(4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235789-85-2 HCAPLUS

CN Acetamide, N-[[(5S)-3-imidazo[2,1-b]benzothiazol-7-yl-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-86-3 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(2-phenylimidazo[2,1-b]benzothiazol-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-87-4 HCAPLUS

CN Imidazo[2,1-b]benzothiazole-2-carboxylic acid, 7-[(5S)-5[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, ethyl ester (9CI) (CA INDEX NAME)

RN 235789-88-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-methylimidazo[2,1-b]benzothiazol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-89-6 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[2-(trifluoromethyl)imidazo[2,1-b]benzothiazol-7-yl]-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-90-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2,3-dimethylimidazo[2,1-b]benzothiazol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Updated Search

RN 235789-91-0 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3-methyl-2-(trifluoromethyl)imidazo[2,1-b]benzothiazol-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-92-1 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(4H-pyrrolo[2,1-c][1,4]benzoxazin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-93-2 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4,4-dimethyl-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-94-3 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4H-imidazo[2,1-c][1,4]benzothiazin-7-y1)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Updated Search

RN 235789-95-4 HCAPLUS

CN Acetamide, N-[[(5S)-3-(5H-imidazo[1,2-a][3,1]benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-96-5 HCAPLUS

CN 4H-Imidazo[5,1-c][1,4]benzoxazine-3-carboxylic acid, 7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, ethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-97-6 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4H-imidazo[5,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235789-98-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(9H-imidazo[1,2-a]indol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235789-99-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(9H-imidazo[1,2-a]indol-7-yl)-2-oxo-5-oxazolidinyl]methyl]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

RN 235790-00-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(5H-imidazo[1,2-a][3,1]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235790-01-9 HCAPLUS

CN Carbamic acid, [7-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-4H-imidazo[2,1-c][1,4]benzoxazin-2-yl]-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235790-02-0 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(4H-pyrazolo[5,1-c][1,4]benzoxazin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235790-03-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4-oxido-5H-imidazo[1,2-a][3,1]benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235790-04-2 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4,4-dioxido-5H-imidazo[1,2-a][3,1]benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235790-17-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(1-bromo-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235790-18-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(1,2-dibromo-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

RN 235790-19-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4,4-difluoro-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235790-26-8 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(4H-tetrazolo[5,1-c][1,4]benzoxazin-7-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 235790-27-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-[2-(hydroxymethyl)-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 235792-17-3 HCAPLUS

CN Acetamide, N-[[(5S)-3-imidazo[1,2-a]quinolin-7-yl-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 235791-13-6P 235791-18-1P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of oxazolidines substituted with bicycles as antimicrobial agents)

RN 235791-13-6 HCAPLUS

CN Acetamide, N-[[(5S)-3-[2-[[[(1,1-dimethylethyl)dimethylsilyl]oxy]methyl]-4H-imidazo[2,1-c][1,4]benzoxazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]-(9CI)
(CA INDEX NAME)

Absolute stereochemistry.

RN 235791-18-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-amino-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

REFERENCE COUNT:

THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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1

ACCESSION NUMBER:

1999:487539 HCAPLUS

DOCUMENT NUMBER:

131:129999

TITLE:

Preparation of oxazolidines substituted with bicycles

as antimicrobial agents

INVENTOR(S):

Bartel, Stephan; Guarnieri, Walter; Haebich, Dieter; Raddatz, Siegfried; Riedl, Bernd; Rosentreter, Ulrich;

Ruppelt, Martin; Stolle, Andreas; Wild, Hanno;

Endermann, Rainer; Kroll, Hein-Peter

PATENT ASSIGNEE(S):

SOURCE:

Bayer A.-G., Germany Ger. Offen., 88 pp.

CODEN: GWXXBX

DOCUMENT TYPE:

Patent German

LANGUAGE:

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FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

		APPLICATION NO.	DATE
WO 9937641	A1 19990729 A1 19990729	DE 1998-19802239 WO 1999-EP96	19990109
DK, EE, ES KE, KG, KP MW, MX, NO	FI, GB, GD, GE, KR, KZ, LC, LK, NZ, PL, PT, RO,	BG, BR, BY, CA, CH, GH, GM, HR, HU, ID, LR, LS, LT, LU, LV, RU, SD, SE, SG, SI, YU, ZW, AM, AZ, BY,	IL, IN, IS, JP, MD, MG, MK, MN, SK, SL, TJ, TM,
FI, FR, GB		UG, ZW, AT, BE, CH, MC, NL, PT, SE, BF, SN, TD, TG	
AU 9926161 EP 1049692	A1 19990809 A1 20001108	AU 1999-26161 EP 1999-906112	19990109
IE, FI		GB, GR, IT, LI, LU,	
JP 2002501065 PRIORITY APPLN. INFO.:	T2 20020115	JP 2000-528563 DE 1998-19802239 WO 1999-EP96	A · 19980122
OTHER SOURCE(S):	MARPAT 131:1299	99	

- * STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY AVAILABLE VIA OFFLINE PRINT *
- AB Title compds. [I; R = Me, Et, COOMe, (CH3)2CH, R1 = H; R-R1 = (CH2)3; R3 = OH, OSO2Me, N3, NH2, NHCOOH2, NHAC, NHCOCH2Br, NHCOOMe, NHCOEt, NHCOCF3, NHCOOBu-t, cyclopropylcarbonylamino, 2-furylcarbonylamino, NHP(OMe)2:O, etc.; X = C:O, CH2, S, S:O, SO2; Y = CH2, CHCH3, NMe, C:O, C6H5CH:C,

C6H5CH2C:, 4-C1C6H4CH:C, 4-MeOC6H4CH:C, etc.; Z = O, CH2], enantiomers, and salts are prepared Thus, the title compound II was prepared from 7-nitro-2H-1, 4-benzoxazin-3-one, C6H5CH2OCONHCl, and (R)-(-)-glycidyl butyrate via reduction cyclization and was tested against Staphylococcus aureus, Mycobacterium smegmatis, and Streptococcus pneumoniae.

IT 233773-77-8P 233773-82-5P 233774-47-5P

233775-01-4P 233775-06-9P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent)

(preparation of oxazolidines substituted with bicycles as antimicrobial agents)

RN 233773-77-8 HCAPLUS

CN 4H-1,4-Benzoxazine-4-carboxylic acid, 7-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,3-dihydro-3-oxo-, methyl ester (9CI) (CA INDEX NAME)

AcNH-CH2

RN 233773-82-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-2H-1,4-benzoxazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-47-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 233775-01-4 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233775-06-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3,4-dihydro-4-methyl-3-oxo-2-(phenylmethylene)-2H-1,4-benzothiazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

IT 233773-46-1P 233773-49-4P 233773-60-9P 233773-61-0P 233773-68-7P 233773-69-8P 233773-83-6P 233773-84-7P 233773-85-8P 233773-92-7P 233773-97-2P 233773-98-3P 233774-06-6P 233774-07-7P 233774-14-6P 233774-15-7P 233774-21-5P 233774-28-2P 233774-29-3P 233774-50-0P 233774-55-5P 233774-56-6P 233774-59-9P 233774-60-2P 233775-02-5P 233775-03-6P 233775-04-7P 233775-05-8P 233775-07-0P 233775-08-1P 233775-09-2P 233775-10-5P 233775-15-0P 233775-16-1P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological

study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

(preparation of oxazolidines substituted with bicycles as antimicrobial agents)

233773-46-1 HCAPLUS RN

Acetamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzoxazin-7-yl)-CN 2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

233773-49-4 HCAPLUS RN

Propanamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzoxazin-7-methyl-3-methyl-CN yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

233773-60-9 HCAPLUS RN

Propanamide, N-[[(5S)-3-(4-ethyl-3,4-dihydro-3-oxo-2H-1,4-benzoxazin-7-yl)-CN 2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Updated Search

233773-61-0 HCAPLUS RN

Acetamide, N-[[(5S)-3-(4-ethyl-3,4-dihydro-3-oxo-2H-1,4-benzoxazin-7-yl)-2-herakide, N-[[(5S)-3-(4-ethyl-3,4-benzoxazin-7-yl)-2-herakide, N-[(5S)-3-(4-ethyl-3-(4-ethyCN oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233773-68-7 HCAPLUS

Propanamide, N-[[(5S)-3-(3,4-dihydro-2,4-dimethyl-3-oxo-2H-1,4-benzoxazin-CN 7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

233773-69-8 HCAPLUS RN

Acetamide, N-[[(5S)-3-(3,4-dihydro-2,4-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-oxo-2H-1,4-benzoxazin-7-dimethyl-3-CN yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233773-83-6 HCAPLUS

Acetamide, N-[(5S)-3-(3,4-dihydro-4-methyl-2H-1,4-benzoxazin-7-yl)-2-oxo-4-methyl-2-oxo-4-methyl-2-benzoxazin-7-yl)-2-oxo-4-methyl-2-benzoxazin-7-yl)-2-oxo-4-methyl-2-benzoxazin-7-yl)-2-oxo-4-methyl-2-benzoxazin-7-yl)-2-oxo-4-methyl-2-benzoxazin-7-yl)-2-oxo-4-methyl-2-benzoxazin-7-yl)-2-oxo-4-methyl-2-benzoxazin-7-yl-CN 5-oxazolidinyl]methyl]-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

•x HCl

RN 233773-84-7 HCAPLUS

CN oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

233773-85-8 HCAPLUS RN

Propanamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-2H-1,4-benzoxazin-7-yl)-2-methyl-2-mCN oxo-5-oxazolidinyl]methyl]-, hydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HC1

RN 233773-92-7 HCAPLUS

CN 4H-1,4-Benzoxazine-4-carboxylic acid, 7-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,3-dihydro-, methyl ester (9CI) (CA INDEX NAME)

RN 233773-97-2 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(2,3,6,7-tetrahydro-3-oxo-5H-pyrido[1,2,3-de]-1,4-benzoxazin-9-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233773-98-3 HCAPLUS

CN Propanamide, N-[[(5S)-2-oxo-3-(2,3,6,7-tetrahydro-3-oxo-5H-pyrido[1,2,3-de]-1,4-benzoxazin-9-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-06-6 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(2,3,6,7-tetrahydro-5H-pyrido[1,2,3-de]-1,4-benzoxazin-9-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-07-7 HCAPLUS

CN Propanamide, N-[[(5S)-2-oxo-3-(2,3,6,7-tetrahydro-5H-pyrido[1,2,3-de]-1,4-benzoxazin-9-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-14-6 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-1-methyl-2-oxo-6-quinolinyl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-15-7 HCAPLUS

CN Propanamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-1-methyl-2-oxo-6-quinolinyl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-21-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(1-ethyl-1,2,3,4-tetrahydro-2-oxo-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-22-6 HCAPLUS

CN Propanamide, N-[[(5S)-3-(1-ethyl-1,2,3,4-tetrahydro-2-oxo-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-28-2 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[1,2,3,4-tetrahydro-1-(1-methylethyl)-2-oxo-6-quinolinyl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

233774-29-3 HCAPLUS RN

CN 6-quinolinyl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

233774-48-6 HCAPLUS RN

Acetamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-1,1-dioxido-2H-1,4-CN benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-50-0 HCAPLUS

Propanamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzothiazin-7-methyl-3-methylCN y1)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 233774-55-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-ethyl-3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-56-6 HCAPLUS

CN Propanamide, N-[[(5S)-3-(2-ethyl-3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233774-59-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-2,4-dimethyl-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

RN 233774-60-2 HCAPLUS

CN Propanamide, N-[[(5S)-3-(3,4-dihydro-2,4-dimethyl-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233775-02-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-(4-ethyl-3,4-dihydro-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233775-03-6 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3,4-dihydro-4-(1-methylethyl)-3-oxo-2H-1,4-benzothiazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 233775-04-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-1,1-dioxido-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233775-05-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-4-methyl-1-oxido-3-oxo-2H-1,4-benzothiazin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233775-07-0 HCAPLUS

CN Acetamide, N-[[(5S)-3-[2-[(4-chlorophenyl)methylene]-3,4-dihydro-4-methyl-3-oxo-2H-1,4-benzothiazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

RN 233775-08-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3,4-dihydro-2-[(4-methoxyphenyl)methylene]-4-methyl-3-oxo-2H-1,4-benzothiazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

RN 233775-09-2 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3,4-dihydro-4-methyl-3-oxo-2-(4-pyridinylmethylene)-2H-1,4-benzothiazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry unknown.

RN 233775-10-5 HCAPLUS

CN Acetamide, N-[[(5S)-3-[3,4-dihydro-4-methyl-3-oxo-2-(phenylmethyl)-2H-1,4-benzothiazin-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 233775-15-0 HCAPLUS
CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-1,3-dimethyl-2-oxo-6-quinazolinyl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 233775-16-1 HCAPLUS
CN Propanamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-1,3-dimethyl-2-oxo-6-quinazolinyl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 31 OF 50

ACCESSION NUMBER:

1999:487536 HCAPLUS

DOCUMENT NUMBER:

131:129985

TITLE:

Oxazolidines substituted by tricyclic indoles

Ruppelt, Martin; Bartel, Stephan; Guarnieri, Walter; Raddatz, Siegfried; Rosentreter, Ulrich; Wild, Hanno; INVENTOR(S):

Endermann, Rainer; Kroll, Hein-Peter

PATENT ASSIGNEE(S):

SOURCE:

Bayer A.-G., Germany Ger. Offen., 40 pp.

CODEN: GWXXBX

DOCUMENT TYPE:

Patent.

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT				KIN		DATE			APPL	ICAT	ION I	NO.		. D.	ATE	
DE 198 WO 993	02235			A1		1999	0729 0729		DE 1	998-	1980	2235		1 1	9980: 9990:	122 109
W:	AL, DK, KE, MW,	AM, EE, KG, MX,	AT, ES, KP, NO,	AU, FI, KR, NZ,	AZ, GB, KZ, PL,	BA, GD, LC, PT,	BB, GE, LK, RO,	BG, GH, LR, RU,	BR, GM, LS, SD,	BY, HR, LT, SE, AM,	CA, HU, LU, SG,	CH, ID, LV, SI,	CN, IL, MD, SK,	CU, IN, MG, SL,	CZ, IS, MK, TJ,	DE, JP, MN, TM,
AU 992	CM, 4206	GM, FR, GA,	GB, GN,	GR, GW, A1	IE, ML,	IT, MR, 1999	LU, NE, 0809	MC, SN,	NL, TD, AU 1	PT, TG .999-	SE, 2420	BF,	ВJ,	CF,	CG, 9990	CI, 109
	AT, IE,	BE, FI	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,
JP 200 PRIORITY AF	PLN.	INFO	.:						DE 1	2000- 1998- 1999-	1980	2235	i	A 1	9990 9980 9990	122

GΙ

$$R_N$$
 N
 O
 CH_2-NH-R^1
 I

AB Approx. 25 antibacterial title compds. such as I (R = benzyl, p-methoxybenzyl, allyl, Bu, cyclohexyl, Et, Me; R1 = Ac, EtCO, CO2Me) were prepared E.g., N-[3-(2-(ethoxycarbonyl)-5-indolylamino)-2-hydroxypropyl]acetamide was cyclized with carbonyldimidazole to give 85% 3-(2-ethoxycarbonyl-5-indolyl)-5-(acetaminomethyl)-2-oxazolidinone. The MIC·of I (R = Bu, R1 = Ac) was 4 μg/mL against Staphylococcus Aureus.

TT 234770-13-9P 234770-20-8P 234770-22-0P 234770-24-2P 234770-26-4P 234770-28-6P 234770-29-7P 234770-35-5P 234770-37-7P 234770-39-9P 234770-40-2P 234770-42-4P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

(preparation and bactericidal activity of oxazolidines substituted by tricyclic indoles)

RN 234770-13-9 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-1H-[1,4]oxazino[4,3-a]indol-8-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 234770-20-8 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1,3-dioxo-2-(2-propenyl)-1H-imidazo[1,5-a]indol-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Acnh-CH₂
$$\stackrel{\circ}{\longrightarrow}$$
 $\stackrel{\circ}{\longrightarrow}$ $\stackrel{\longrightarrow$

RN 234770-22-0 HCAPLUS

CN Acetamide, N-[[3-(2-butyl-2,3-dihydro-1,3-dioxo-1H-imidazo[1,5-a]indol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 234770-24-2 HCAPLUS

CN Acetamide, N-[[3-(2-ethyl-2,3-dihydro-1-oxo-3-thioxo-1H-imidazo[1,5-a]indol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 234770-26-4 HCAPLUS

CN Acetamide, N-[[(5S)-3-(3,4-dihydro-1-oxo-1H-[1,4]oxazino[4,3-a]indol-8-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 234770-28-6 HCAPLUS

CN Propanamide, N-[[(5S)-3-(3,4-dihydro-1-oxo-1H-[1,4]oxazino[4,3-a]indol-8-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 234770-29-7 HCAPLUS

Acetamide, N-[[3-(2,3-dihydro-2-methyl-3-oxo-1H-imidazo[1,5-a]indol-7-yl)-CN 2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

234770-35-5 HCAPLUS RN

Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-2-methyl-1-oxopyrazino[1,2-tetrahydro-2-methyl-1-oxopyraziCN a]indol-8-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

234770-37-7 HCAPLUS RN

Propanamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-2-methyl-1-CN oxopyrazino[1,2-a]indol-8-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

234770-39-9 HCAPLUS RN

Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-2-methyl-3-oxopyrazino[1,2-methyl-3-oxopyrazino]]CN a]indol-8-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 234770-40-2 HCAPLUS
CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-2-methyl-1,4-dioxopyrazino[1,2-a]indol-8-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 234770-42-4 HCAPLUS
CN Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-1-oxopyrazino[1,2-a]indol-8-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 234770-47-9P
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
 (Reactant or reagent)

(preparation and bactericidal activity of oxazolidines substituted by tricyclic indoles)

RN 234770-47-9 HCAPLUS

CN 1H-Indole-2-carboxylic acid, 5-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, ethyl ester (9CI) (CA INDEX NAME)

AcNH-CH2

IT 234770-15-1P 234770-17-3P 234770-23-1P

234770-25-3P 234770-41-3P

RL: SPN (Synthetic preparation); PREP (Preparation)

(preparation and bactericidal activity of oxazolidines substituted by tricyclic indoles)

RN 234770-15-1 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1,3-dioxo-2-(phenylmethyl)-1H-imidazo[1,5-a]indol-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 234770-17-3 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-[(4-methoxyphenyl)methyl]-1,3-dioxo-1H-imidazo[1,5-a]indol-7-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 234770-23-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-(2-cyclohexyl-2,3-dihydro-1,3-dioxo-1H-imidazo[1,5-a]indol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 234770-25-3 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2-methyl-1,3-dioxo-1H-imidazo[1,5-a]indol-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

234770-41-3 HCAPLUS RN

Acetamide, N-[[(5S)-2-oxo-3-(1,2,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,3,4-tetrahydro-2-methyl-1,4-tetCN trioxopyrazino[1,2-a]indol-8-yl)-5-oxazolidinyl]methylj- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

L23 ANSWER 32 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1999:487281 HCAPLUS

DOCUMENT NUMBER:

131:116228

TITLE:

Preparation of oxazolidinones as bactericides

INVENTOR(S):

Gordeev, Mikhail F.; Luehr, Gary W.; Patel, Dinesh V.;

Ni, Zhi-Jie; Gordon, Eric

PATENT ASSIGNEE(S):

SOURCE:

Versicor, Inc., USA PCT Int. Appl., 229 pp.

CODEN: PIXXD2 Patent

DOCUMENT TYPE:

LANGUAGE:

English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PAT	ENT	NO.			KINI) .	DATE			APPL	ICAT:	ION I	NO.		D2	ATE	
WO 9937630 A1 1999072						0729	WO 1999-US1318						19990122				
					AU,												
		DK,	EE,	ES,	FI,	GB,	GD,	GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,
					KR,												
		MW,	MX,	NO,	ΝZ,	PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	ТJ,	TM,
		TR,	TT,	UA,	UG,	US,	UZ,	VN,	YU,	zw							
	RW:	GH,	GM,	ΚE,	LS,	MW,	SD,	SZ,	UG,	ZW,	ΑT,	BE,	CH,	CY,	DE,	DK,	ES,
		FI,	FR,	GB,	GR,	ΙE,	ΙT,	LU,	MC,	NL,	PT,	SE,	BF,	ВJ,	CF,	CG,	CI,
					GW,												
CA	2318	969			AA		1999	0729		CA 1	999-:	2318	969		1	9990	122
ΑU	9924	644			A1		1999	0809		AU 1	9.99-	2464	4		1	9990	122
ΑU	7641	84			В2		2003	0814									
ΕP	1049	682			A1		2000	1108		EP 1	999-	9041	93		1	9990	122
	R:	ΑT,	BE,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,
		ΙE,	SI,	LT,	LV,	FI,	RO										

JP 2002501059 BR 9907183 NZ 505902 PRIORITY APPLN. INFO.:	T2 A A	20020115 20030610 20030829	BR NZ US	2000-528553 1999-7183 1999-505902 1998-12535 1998-86702 1999-US1318	A	19990122 19990122 19990122 19980123 19980528 19990122
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OTHER SOURCE(S):

MARPAT 131:116228

GI

ΙT

Title compds. [e.g., I; R = H; R1 -SR11, CONR7R8, etc.; R7,R8. R11 = H, alkyl, (hetero)aryl, etc.] were prepared Thus, 3,4-F(Me3CO2C)C6H3NHCO2CH2Ph (preparation given) was cyclocondensed with (R)-glycidyl butyrate and the product converted in several steps to I (R = resin, R1 = CO2C6F5) which was amidated by morpholine to give, after resin cleavage, I (R = H, R1 = CONHR8, R8 = morpholino). Data for biol. activity of I were given.

RN 232950-89-9 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-[1-(triphenylmethyl)-1H-indazol-5-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT:

14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L23 ANSWER 33 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1999:421579 HCAPLUS

DOCUMENT NUMBER:

131:78435

TITLE:

Method and device for in-situ formulation of a medicinal solution for parenteral application Kuehn, Bernd; Wiessmeier, Georg; Rupp, Roland;

INVENTOR(S):

Krumbach, Bernd; Weismantel, Lothar; Herrmann, Erhard;

Klein, Juergen

PATENT ASSIGNEE(S):

Bayer Aktiengesellschaft, Germany

SOURCE:

PCT Int. Appl., 47 pp.

CODEN: PIXXD2

DOCUMENT TYPE:

Patent German

LANGUAGE:

. 1

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PAS	CENT 1	NO.			KIN)	DATE			APP.	LICAT	ION I	NO.		D.	ATE	
WO	9932	 175			A1	-	1999	0701	,	WO	1998-	EP80	14		1	9981	209
	W:	AL.	AM.	AT,	AU,	AZ,	BA,	BB,	BG,	BR	, BY,	CA,	CH,	CN,	CU,	CZ,	DE,
		DK.	EE,	ES.	FI,	GB,	GD,	GE,	GH,	GM	, HR,	HU,	ID,	IL,	IN,	IS,	JP,
		KE.	KG.	KP.	KR,	KZ,	LC,	LK,	LR,	LS	, LT,	LU,	LV,	MD,	MG,	MK,	MN,
		MW,	MX,	NO,	NZ,	PL,	PT,	RO,	RU,	SD	, SE,	SG,	SI,	SK,	SL,	ТJ,	TM,
		TR,	TT,	UA,	UG,	US,	UZ,	VN,	YU,	ZW	, AM,	ΑZ,	BY,	KG,	ΚZ,	MD,	RU,
		TJ,	TM	-													
	RW:	GH,	GM,	KE,	LS,	MW,	SD,	SZ,	UG,	ZW	, AT,	BE,	CH,	CY,	DE,	DK,	ES,
		FI,	FR,	GB,	GR,	ΙE,	IT,	LU,	MC,	NL	, PT,	SE,	BF,	ВJ,	CF,	CG,	CI,
		CM,	GA,	GN,	GW,	ML,	MR,	ΝE,	SN,	TD	, TG						
DE	1975						1999	0701		DE	1997-				1	9971	
AU	9922	694			A1		1999	0712		ΑU	1999-	2269	4			9981	
ĖP	1045	709			A1		2000	1025		EΡ	1998-	9662	73		1	9981	209
EP	1045	709			В1		2003	1203									
	R:	DE,	ES,	FR,	GB,	ΙT											
JP	2001				Т2		2001	1218			2000-					9981	
ES	2212	390			т3		2004	0716			1998-				_	9981	
US	6540	715			В1		2003	0401			2000-				_	0000	
PRIORIT	Y APP	LN.	INFO	.:							1997-					9971	
										WO	1998-	EP80	14		W 1	9981	209

AB In the title method, ≥2 dosed partial flows are continually incorporated by a mixer into a total volume flow containing active substances

form a medicinal infusion solution which is not in thermodn. equilibrium; the resulting total volume flow after mixing is 0.2-500 mL/h, preferably 5-500 mL/h. One of the partial flows may be an active substance concentrate in an organic or aqueous-organic solvent, whereas the other partial flow comprises an aqueous

or aqueous-organic diluent; the nonequil. state may comprise a supersatd. solution

The mixer is preferably one free of dead space, such as an orifice-type blender. An active substance administration set comprises the mixer, infusion tubing, storage containers for the active substance concentrate and diluent, and packing materials. Thus, a 2% concentrate of oxazolidinone antibiotic Bay 17-1648 in glycofurol was delivered with a perfusion pump to a mixer for dilution 1:13 with water delivered with an infusion pump, to produce a total volume flow of 430 mL effluent/h containing 7% glycofurol and 140 mg Bay 17-1648/100 mL (.apprx.7 times the concentration for saturation);

this

to

solution was stable against crystallization of Bay 17-1648 for ≥ 1 min, and glycofurol at this concentration was well tolerated. Administration of a daily dose of 1000 mg Bay 17-1648 required an infusion volume of 717 mL, which would require .apprx.100 min for infusion.

IT 184157-64-0, Bay 17-1648

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(method and device for in-situ formulation of medicinal solution for parenteral application)

RN 184157-64-0 HCAPLUS

CN Propanamide, N-[[(5S)-3-(2,3-dihydro-3-methyl-2-oxo-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT:

THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 34 OF 50

5

ACCESSION NUMBER:

1999:77555 HCAPLUS

DOCUMENT NUMBER:

130:139335

TITLE:

Preparation of tricyclically substituted

oxazolidinones as bactericides

INVENTOR(S):

Bartel, Stephan; Guarnieri, Walter; Riedl, Bernd; Habich, Dieter; Stolle, Andreas; Ruppelt, Martin; Raddatz, Siegfried; Rosentreter, Ulrich; Wild, Hanno;

Endermann, Rainer; Kroll, Hein-peter

PATENT ASSIGNEE(S):

Bayer Aktiengesellschaft, Germany; et al.

SOURCE:

PCT Int. Appl., 98 pp. CODEN: PIXXD2

DOCUMENT TYPE:

Patent

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT	NO.		KIND	DATE			APPL	ICAT	ION N	10.		D2	ATE	-
WO 9903	 846		A1	1999	0128		WO 1	998-1	EP425	52		1	9980	708
W:	AL, AM,	AT,	AU, F	AZ, BA,	BB,	BG,	BR,	BY,	CA,	CH,	CN,	CU,	CZ,	DE,
	DK, EE,	ES.	FI. C	B, GE,	GH,	GM,	HR,	HU,	ID,	IL,	IS,	JP,	KE,	KG,
	KP, KR,	KZ.	LC, I	K, LR,	LS,	LT,	LU,	LV,	MD,	MG,	MK,	MN,	MW,	MX,
	NO, NZ,	PL,	PT, F	RO, RU,	SD,	SE,	SG,	SI,	SK,	SL,	ТJ,	TM,	TR,	TT,
	UA, UG,	US.	UZ. \	/N, YU,	ZW,	AM,	AZ,	BY,	KG,	ΚZ,	MD,	RU,	ТJ,	TM
RW:	GH, GM,	KE.	LS. N	W, SD.	SZ,	UG,	ZW,	AT,	BE,	CH,	CY,	DE,	DK,	ES,
••	FI, FR,	GB.	GR.	E. IT.	LU,	MC,	NL,	PT,	SE,	BF,	ВJ,	CF,	CG,	CI,
	CM, GA,			1L, MR,										
DE 1973		•	•	1999			DE 1		19730	0847		1	9970	718
AU 9884				1999			AU 1	998-	84417	7		1	9980	708
	360		A	1999	0127		ZA 1	998-	6360			1	9980	717
PRIORITY APP							DE 1	997-	19730	0847		A 1	9970	718
11110111111111							wo 1	998-	EP425	52	1	W 1	9980	708
OTHER SOURCE	(S):		MARPA	AT 130:	1393	35								

GΙ

Title compds. [I; R1= N3, OH, OMe, OSO2Me, NH2, NHCOCH3, etc.; E = O, S, CO, SO, SO2, NC2H5, etc.; A, A1, A2, A3 are independently CH, N, with no more than one N; L and M are independently H, OH, CO, CN, NO2, CHO, etc.; dotted bonds = one single bond to I and the other single bond to a H] are prepared as antibacterial medicaments. Thus, compound II was prepared from cycloaddn. of 2-benzyloxycarbonylaminofluorene and (R)-2,3-epoxypropyl butanoate in the presence of Bu lithium in hexane.

IT 220059-31-4P 220059-56-3P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent) (preparation of tricyclically substituted oxazolidinones as bactericides)

RN 220059-31-4 HCAPLUS
CN Acetamide, N-[[(5S)-3-(2-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl](9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-56-3 HCAPLUS
CN Acetamide, N-[[(5S)-2-oxo-3-[7-(3-pyridinyl)-9H-fluoren-2-yl]-5oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

220058-98-0P 220059-01-8P 220059-05-2P IT 220059-07-4P 220059-08-5P 220059-09-6P 220059-12-1P 220059-13-2P 220059-16-5P 220059-17-6P 220059-18-7P 220059-19-8P 220059-25-6P 220059-26-7P 220059-29-0P 220059-33-6P 220059-34-7P 220059-35-8P 220059-36-9P 220059-37-0P 220059-38-1P 220059-39-2P 220059-42-7P 220059-43-8P 220059-46-1P 220059-47-2P 220059-50-7P 220059-51-8P 220059-55-2P 220059-59-6P 220059-60-9P 220059-61-0P 220059-62-1P 220059-64-3P RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation) (preparation of tricyclically substituted oxazolidinones as bactericides) 220058-98-0 HCAPLUS RN Acetamide, N-[[(5S)-3-(9H-fluoren-2-yl)-2-oxo-5-oxazolidinyl]methyl]-CN (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-01-8 HCAPLUS CN Propanamide, N-[[(5S)-3-(9H-fluoren-2-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-05-2 HCAPLUS
CN Acetamide, N-[[(5S)-3-(9-ethyl-9H-carbazol-3-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-07-4 HCAPLUS

CN Propanamide, N-[[(5S)-3-(9-ethyl-9H-carbazol-3-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-08-5 HCAPLUS.

CN Acetamide, N-[[(5S)-3-(7-acetyl-9H-fluoren-2-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-09-6 HCAPLUS

CN Propanamide, N-[[(5S)-3-(7-acetyl-9H-fluoren-2-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

220059-12-1 HCAPLUS RN

Acetamide, N-[[(5S)-3-(3-dibenzofurany1)-2-oxo-5-oxazolidiny1]methy1]-CN (CA INDEX NAME)

Absolute stereochemistry.

220059-13-2 HCAPLUS RN

Propanamide, N-[[(5S)-3-(3-dibenzofuranyl)-2-oxo-5-oxazolidinyl]methyl]-1CN (9CI) (CA INDEX NAME)

Absolute stereochemistry.

220059-16-5 HCAPLUS RN

Acetamide, N-[[(5S)-2-oxo-3-[8-(1E)-1-pentenyl-3-dibenzofuranyl]-5-1-pentenyl-3-dibenzofuranyl]CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry as shown.

RN 220059-17-6 HCAPLUS CN Propanamide, N-[[(5S)-2-oxo-3-[8-(1E)-1-pentenyl-3-dibenzofuranyl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry as shown.

RN 220059-18-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(8-acetyl-3-dibenzofuranyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-19-8_ HCAPLUS

CN Propanamide, N-[[(5S)-3-(8-acetyl-3-dibenzofuranyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-25-6 HCAPLUS

CN Acetamide, N-[[(5S)-2-oxo-3-(9-oxo-9H-fluoren-2-yl)-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

RN 220059-26-7 HCAPLUS
CN Propanamide, N-[[(5S)-2-oxo-3-(9-oxo-9H-fluoren-2-yl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-29-0 HCAPLUS
CN Propanamide, N-[[(5S)-3-(2-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl](9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-33-6 HCAPLUS CN Acetamide, N-[[(5S)-3-(9H-carbazol-3-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-34-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-[7-(dimethylamino)-9H-fluoren-2-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-35-8 HCAPLUS

CN Acetamide, N-[[(5S)-3-(7-cyano-9H-fluoren-2-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-36-9 HCAPLUS

CN Propanamide, N-[[(5S)-3-[7-(dimethylamino)-9H-fluoren-2-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-37-0 HCAPLUS

CN Acetamide, N-[[(5S)-3-(7-bromo-9H-fluoren-2-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

220059-38-1 HCAPLUS RN Propanamide, N-[[(5S)-3-(3-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl]-CN (9CI) (CA INDEX NAME)

Absolute stereochemistry.

220059-39-2 HCAPLUS RN

Acetamide, N-[[(5S)-3-(3-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl]-CN (CA INDEX NAME)

Absolute stereochemistry.

220059-42-7 HCAPLUS RN

Acetamide, N-[[(5S)-3-(9-methyl-9H-carbazol-2-yl)-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-43-8 HCAPLUS

Propanamide, N-[[(5S)-3-(9-methyl-9H-carbazol-2-yl)-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 220059-46-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-(5H-indeno[1,2-b]pyridin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-47-2 HCAPLUS

CN Propanamide, N-[[(5S)-3-(5H-indeno[1,2-b]pyridin-7-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-50-7 HCAPLUS

CN Acetamide, N-[[(5S)-3-(5,5-dioxido-2-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-51-8 HCAPLUS
CN Acetamide, N-[[(5S)-3-(5-oxido-2-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-55-2 HCAPLUS CN Propanamide, N-[[(5S)-2-oxo-3-[7-(3-pyridinyl)-9H-fluoren-2-yl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-59-6 HCAPLUS CN Acetamide, N-[[(5S)-2-oxo-3-[7-(3-pyridinyl)-9H-fluoren-2-yl]-5-oxazolidinyl]methyl]-, monohydrochloride (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

RN 220059-60-9 HCAPLUS CN Acetamide, N-[[(5S)-3-(5,5-dioxido-3-dibenzothienyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA_INDEX_NAME)

Absolute stereochemistry.

RN 220059-61-0 HCAPLUS

CN Acetamide, N-[[(5S)-3-benzofuro[2,3-b]pyridin-7-yl-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-62-1 HCAPLUS

CN Acetamide, N-[[(5S)-3-benzofuro[3,2-c]pyridin-7-yl-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 220059-64-3 HCAPLUS

CN Acetamide, N-[[(5S)-3-benzofuro[2,3-c]pyridin-7-yl-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

REFERENCE COUNT:

THERE ARE 6 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L23 ANSWER 35 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

6

ACCESSION NUMBER: 1997:552637 HCAPLUS

DOCUMENT NUMBER: 127:149139

TITLE: Preparation of bactericidal pyridothienyl- and

pyridofuryloxazolidinones

INVENTOR(S): Riedl, Bernd; Haebich, Dieter; Stolle, Andreas;

Ruppelt, Martin; Bartel, Stefan; Guarnieri, Walter;

Endermann, Rainer; Kroll, Hein-Peter

PATENT ASSIGNEE(S): Bayer A.-G., Germany SOURCE: Ger. Offen., 28 pp.

CODEN: GWXXBX

DOCUMENT TYPE: Patent

LANGUAGE: German FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PA	TENT	NO.	•		KIND)	DATE			APE	PLI	CAT	ION	NO.		D	ATE	
DE	1960	 1264			A1		1997	0717		DE	19	96-	1960	1264		1	9960	116
	7852				A2		1997	0723		ΕP	19	97-	1000	25		1	9970	103
	7852				A 3		1999	0210										
		AT,	BE,	CH,	DE,	DK	, ES,	FI,	FR	, GE	3,	GR,	ΙE,	ΙΤ,	LI,	LU,	MC,	NL,
		PT,	SE															
AU	9710	098			A1		1997	0724		ΑU	19	97-	1009	8		_	9970	
US	5827	857			Α		1998	1027					7810			_	9970	
	2194				AA		1997	0717		CA	19	97-	2194	938		_	9970	
JP	0919	4482			A2		1997	0729	•	JΡ	19	97-	1755	9		1	9970	114
NO	9700	175			Α		1997	0717		NO	19	97-	175			_	9970	
7.A	9700	303			Α		1997	0717		ZA	19	97-	303			1	.9970	115
CN	1161	968			Α		1997	1015		CN	19	97-	1018	06		1	.9970	116
	9700	-			A ·		1998	0901		BR	19	97-	702			1	.9970	116
PRIORIT			INFO	. :						DE	19	96-	1960	1264		A 1	.9960	116
OTHER SO					MARE	TAS	127:	1491	39									
O 1																		

$$RN \longrightarrow O$$
 CH_2R^1 I CH_2R^2 II

Title compds. I [R = pyridothienyl, pyridofuryl; R1 = N3, (un)substituted OH, NH2] were prepared Thus, 2-chloro-6-methylpyridine-3-carebonitrile was treated with HSCH2CO2Me to give Me 3-amino-6-methylthieno[2,3-b]pyridine-2-carboxylate, which was deaminated, and converted to the 2-butoxycarbonylamino derivative via the acid azide. 2-Butoxycarbonylamino-6-methylthieno[2,3-b]pyridine was cyclized with (R)-glycidyl butyrate to give the oxazolidinone II [R2 = OH]. This was converted to II [R2 = NHCSMe] which had min. inhibitory concns. against several staphylococcus strains of 2 μ g/mL.

IT 193400-72-5P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of bactericidal pyridothienyl- and pyridofuryloxazolidinones)

RN 193400-72-5 HCAPLUS

CN Acetamide, N-[[3-(5-methylthieno[2,3-b]pyridin-2-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Absolute stereochemistry.

RN 193400-79-2 HCAPLUS
CN Propanamide, N-[[3-(6-methylthieno[2,3-b]pyridin-2-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193400-82-7 HCAPLUS
CN Acetamide, N-[[3-(6-methylthieno[2,3-b]pyridin-2-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 36 OF 50

ACCESSION NUMBER:

1997:552636 HCAPLUS

DOCUMENT NUMBER:

127:149138

TITLE:

Preparation of antimicrobial dihydroquinolinyloxazolidinones

INVENTOR(S):

Haebich, Dieter; Stolle, Andreas; Riedl, Bernd; Ruppelt, Martin; Bartel, Stefan; Guarnieri, Walter; Endermann, Rainer; Kroll, Hein-Peter

PATENT ASSIGNEE(S):

Bayer A.-G., Germany

SOURCE:

Ger. Offen., 29 pp. CODEN: GWXXBX

DOCUMENT TYPE:

Patent

LANGUAGE:

German

1

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 19601265 EP 785197	A1 A2	19970717 19970723	DE 1996-19601265 EP 1997-100026	. 19960116 19970103
EP 785197	A 3	19990210		
R: AT, BE, CH,	DE, DK	, ES, FI,	FR, GB, GR, IE, IT, LI	, LU, MC, NL,
PT, SE				10070100
US 5861413	Α	19990119	US 1997-781002	19970109
AU 711924	B2	19991021	AU 1997-10097	19970109
CA 2194945	AA	19970717	CA 1997-2194945	19970113
JP 09194478	A2	19970729	JP 1997-15889	19970113
NO 9700174	A	19970717	NO 1997-174	19970115
ZA 9700302	A	19970717	ZA 1997-302	19970115
BR 9700688	A	19980901	BR 1997-688	19970115
ни 9700126	A2	19981228	HU 1997-126	19970115
CN 1163892	. A	19971105	CN 1997-102291	19970116
PRIORITY APPLN. INFO.:			DE 1996-19601265	A 19960116
OTHER SOURCE(S):	MARPAT	127:14913	38	
GI				

AB Title compds. were prepared Thus, 6-aminoquinoline was N-benzyloxycarbonylated and treated with (R)-glycidyl butyrate to give (5R)-3-(6-quinolinyl)-5-hydroxymethyl-2-oxazolidinone which was converted to the acetylaminomethyl deriv.in 4 steps and then to the dihydroquinoline via the N-oxide, followed by methylation to give the title compound I. I had min. inhibitory concns. against Staphylococcus 133 and Mycobacterium smegmatis of 2 μg/mL.

IT 193359-44-3P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(preparation of antimicrobial dihydroquinolinyloxazolidinones)

RN 193359-44-3 HCAPLUS

CN Acetamide, N-[[3-[1-(cyanomethyl)-1,2-dihydro-2-oxo-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Absolute stereochemistry.

RN 193359-40-9 HCAPLUS
CN Acetamide, N-[[3-(1-ethyl-1,2-dihydro-2-oxo-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

193359-42-1 HCAPLUS RN

CN 5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-46-5 HCAPLUS

Acetamide, N-[[3-[1,2-dihydro-1-(2-hydroxyethyl)-2-oxo-6-quinolinyl]-2-oxo-6-quinolinyl]CN 5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

193359-56-7 HCAPLUS RN

Acetamide, N-[[3-[1,2-dihydro-1-(methylsulfonyl)-2-oxo-6-quinolinyl]-2-oxo-6-quinolinyl]CN 5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

IT 175391-92-1P 175392-27-5P 193359-34-1P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation of antimicrobial dihydroquinolinyloxazolidinones)

RN 175391-92-1 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(6-quinolinyl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-27-5 HCAPLUS

CN Acetamide, N-[[3-(1-oxido-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-34-1 HCAPLUS

CN Acetamide, N-[[3-(1,2-dihydro-2-oxo-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

IT 193359-48-7P 193359-50-1P 193359-52-3P

193359-55-6P 193359-58-9P 193359-61-4P

RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of antimicrobial dihydroquinolinyloxazolidinones)

RN 193359-48-7 HCAPLUS

CN Acetamide, N-[[3-[1,2-dihydro-2-oxo-1-(phenylmethyl)-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-50-1 HCAPLUS

CN 1(2H)-Quinolineacetamide, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-oxo-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-52-3 HCAPLUS

CN Acetamide, N-[[3-[1,2-dihydro-1-(hydroxymethyl)-2-oxo-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-55-6 HCAPLUS

CN Acetamide, N-[[3-[1-[(dimethylamino)methyl]-1,2-dihydro-2-oxo-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-58-9 HCAPLUS

CN Acetamide, N-[[3-[1-(3-chlorobenzoyl)-1,2-dihydro-2-oxo-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 193359-61-4 HCAPLUS

CN 1(2H)-Quinolinecarboxamide, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-N-methyl-N-[(methylamino)carbonyl]-2-oxo-, (S)- (9CI) (CFINDEX NAME)

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 37 OF 50

ACCESSION NUMBER:

1997:453982 HCAPLUS

DOCUMENT NUMBER:

127:81464

TITLE:

Preparation of oxazolidinone antibacterial agents with

tricyclic substituents

INVENTOR(S):

Thomas, Richard C.; Cleek, Gary J.; Hutchinson, Douglas K.; Yamada, Hiroyoshi Pharmacia & Upjohn Company, USA

PATENT ASSIGNEE(S): SOURCE:

PCT Int. Appl., 41 pp.

CODEN: PIXXD2

DOCUMENT TYPE:

Patent

LANGUAGE:

English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PA	PATENT NO.				KIND DATE		APPLICATION NO.					DATE						
WO							WO 1996-US17120				19961105							
											BY,							
											JP,							
											MN,							
											TR,							
								RU,										
	RW:										DE,	DK,	ES,	FI,	FR,	GB,	GR,	
											CF,							
		MR,	NE,	SN,	TD,	ΤG												
ZA	9608	661			Α		1998	0414		ZA 1	996-	8661			1	9961	014	
AU	9676	651			A1		1997	0611		AU 1	.996-	7665	1		1	9961	105	
EP	8748	52			A1		1998	1104		EP 1	.996-	9394	96		1	9961	105	•
EP	8748																	
	R:	ΑT,	BE,	CH,	DE,	DK,	ES,	FR,	GB,	GR,	IT,	LI,	LU,	NL,	SE,	MC,	PT,	
		ΙE,	SI,	LT,	LV,	FΙ,	RO											
JP	2000	5004	89		Т2		2000	0118		JP 1	.997-:	5197	26		1	9961	105	
ΑT	2631	72			E		2004	0415		AT 1	.996-	9394	96		1	9961	105	
PT	2000 2631 8748 2217	52			${f T}$		2004	0831		PT 1	996-	9394	96		1	9961	105	
ES	2217	329			т3		2004	1101		ES 1	.996-	9394	96		1	9961	105	
TW	4266	83			В		2001	0321		TW 1	.996-	8511	4048		1	9961	115	
	5922				Α		1999	07.13			997-							
US	5955	460			Α		1999	0921			999-					9990		
CORIT	Y APP	LN.	INFO	.:							995-							
											996-	_				9961		
											996-							
										US 1	1997-	8504	24		A3 1	9970	502	
HER S	OURCE	(S):			MAR	PAT	127:	81464	1									

OTHER SOURCE(S): MARPAT 127:81464

AB The title compds. [I; R = H, alkyl, cycloalkyl, alkoxy, (un) substituted NH2; X = (un) substituted NH, (un) substituted CH2, S, SO, SO2, O; Y = H, halogen; Z = (un) substituted NH, S, SO, SO2, O], which are effective antibiotics against a number of gram-pos. aerobic bacteria as well as anaerobic organisms and acid-fast organisms, are prepared Thus, phenylmethyl 8-[5-(S)-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydropyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylate was prepared in 9 steps from 2-pyrazinecarboxylic acid and demonstrated a MIC against S. aureus UC Number 9213 of 8 μg/mL.

IT 191738-68-8P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(preparation of oxazolidinone antibacterial agents with tricyclic substituents)

RN 191738-68-8 HCAPLUS

CN Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid, 8-[(5S)-5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-, phenylmethyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191738-76-8 HCAPLUS
CN Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid,
8-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-,
methyl ester, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191738-81-5 HCAPLUS
CN Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(hydroxyacetyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

RN 191738-85-9 HCAPLUS
CN Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(5-isoxazolylcarbonyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191738-89-3 HCAPLUS
CN Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(1H-indol-2-ylcarbonyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191738-94-0 HCAPLUS CN Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid,

8-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-, phenylmethyl ester, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191738-96-2 HCAPLUS

CN Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid, 8-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-, phenylmethyl ester, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191738-98-4 HCAPLUS

CN Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid, 8-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-, methyl ester, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

191739-01-2 HCAPLUS RN Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid, CN

8-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-, methyl ester, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-03-4 HCAPLUS

Acetamide, N-[[3-(1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-hexahydropyrazin-8-hexahydCN yl)-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

191739-05-6 HCAPLUS RN

Acetamide, N-[[3-(1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-CN y1)-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

RN 191739-07-8 HCAPLUS

CN Acetamide, N-[[3-(1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl)-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-09-0 HCAPLUS

CN Acetamide, N-[[3-[3-(2-fluoroethyl)-1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-11-4 HCAPLUS

CN Acetamide, N-[[3-[3-(cyanomethyl)-1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

RN 191739-13-6 HCAPLUS
CN Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(methylsulfonyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-15-8 HCAPLUS
CN Acetamide, N-[[3-(3-formyl-1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl)-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CAINDEX NAME)

Absolute stereochemistry.

RN 191739-17-0 HCAPLUS
CN Acetamide, N-[[3-(3-acetyl-1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl)-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CAINDEX NAME)

RN 191739-19-2 HCAPLUS CN Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(methoxyacetyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA

Absolute stereochemistry.

INDEX NAME)

RN 191739-21-6 HCAPLUS

CN Acetamide, N-[[3-[3-[(acetyloxy)acetyl]-1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-23-8 HCAPLUS

CN Acetamide, N-[[3-[3-(dichloroacetyl)-1,2,3,4,4a,5-hexahydropyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-25-0 HCAPLUS

Pyrazino[2,1-c][1,4]benzoxazine-3(4H)-carboxylic acid, 8-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1,2,4a,5-tetrahydro-, 1,1-dimethylethyl ester, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 191739-27-2 HCAPLUS

CN Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(3-hydroxy-1-oxopropyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

191739-29-4 HCAPLUS RN

Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(3-methoxy-1-CN oxopropyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

191739-31-8 HCAPLUS RN

Acetamide, N-[[3-[3-(1,4-dioxopentyl)-1,2,3,4,4a,5-hexahydropyrazino[2,1-dioxopentyl]]CN c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

191739-33-0 HCAPLUS RN

Acetamide, N-[[3-[1,2,3,4,4a,5-hexahydro-3-(5-nitro-2-CN thiazolyl)pyrazino[2,1-c][1,4]benzoxazin-8-yl]-2-oxo-5-oxazolidinyl]methyl]-, [8(S)]- (9CI) (CA INDEX NAME)

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 38 OF 50

ACCESSION NUMBER:

1996:753781 HCAPLUS

DOCUMENT NUMBER:

126:18862

TITLE:

Preparation of N-(oxobenzoxazol-6-yl)oxazolidinones

and analogs as antibacterial agents

INVENTOR(S):

Stolle, Andreas; Haebich, Dieter; Bartel, Stephan; Riedl, Bernd; Ruppelt, Martin; Wild, Hanno; Endermann,

Rainer; Bremm, Klaus-Dieter; Kroll, Hein-Peter; et al.

PATENT ASSIGNEE(S):

Bayer A.-G., Germany Eur. Pat. Appl., 117 pp.

SOURCE:

CODEN: EPXXDW

DOCUMENT TYPE:

Patent

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE		
EP 738726 EP 738726	A1 B1	19961023 20010926	EP 1996-105539	19960409		
R: AT, BE, CH, PT, SE			R, GB, GR, IE, IT, LI,	LU, MC, NL,		
DE 19544106	A1	19961024	DE 1995-19544106	19951127		
AT 206120	E	20011015	AT 1996-105539	19960409		
ES 2164182	т3	20020216	ES 1996-105539	19960409		
US 6069160 '	Α.	20000530	US 1996-631516	19960412		
JP 08301869	A2	19961119	JP 1996-117117	19960416		
AU 9650735	A1	19961031	AU 1996-50735	19960417		
AU 705071	В2	19990513				
ни 9601001	A2	19980428	ни 1996-1001	19960417		
CA 2174473	AA	19961022	CA 1996-2174473	19960418		
NO 9601559	A	19961022	NO 1996-1559	19960419		
ZA 9603138	A	19961104	ZA 1996-3138	19960419		
CN 1138582	A	19961225	CN 1996-106152	19960419		
BR 9602016	A	19980407	BR 1996-2016	19960422		
CN 1161336	A	19971008	CN 1997-102064	19970118		
PRIORITY APPLN. INFO.:	••			A 19950421		
INIONIII MIIDN. INIO				A 19951127		
OTHER COURCE (C).	маррат	126.18862				

OTHER SOURCE(S):

MARPAT 126:18862

GΙ

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Title compds. [I; R = e.g., oxobenzoxazol-6-yl, etc.; R1 = N3,
AB
     (protected) hydroxy, acyloxy, alkylsulfonyloxy, NR4R5, etc.; R4,R5 = H,
     alkyl, Ph, etc.] were prepared Thus, 6-benzyloxycarbonylamino-3-methyl-2-
     benzothiazolinone (preparation given) was cyclocondensed with (R)-glycidyl
     butyrate to give title compound II. Data for antibacterial activity of
     selected I were given.
     184156-99-8P 184157-00-4P 184157-01-5P
IT
     184157-02-6P 184157-03-7P 184157-06-0P
     184157-08-2P 184157-10-6P 184157-12-8P
     184157-14-0P 184157-17-3P 184157-19-5P
     184157-21-9P 184157-23-1P 184157-26-4P
     184157-29-7P 184157-32-2P 184157-35-5P
     184157-38-8P 184157-41-3P 184157-44-6P
     184157-45-7P 184157-47-9P 184157-49-1P
     184157-50-4P 184157-51-5P 184157-52-6P
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     184157-56-0P 184157-57-1P 184157-58-2P
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     184157-96-8P 184157-97-9P 184157-98-0P
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     184158-02-9P 184158-03-0P 184158-04-1P
     184158-05-2P 184158-06-3P 184158-07-4P
     184158-08-5P 184158-09-6P 184158-23-4P
     184158-29-0P 184158-31-4P 184158-39-2P
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     184158-68-7P 184158-69-8P 184158-70-1P
     184158-71-2P 184158-72-3P 184158-73-4P
     184158-78-9P 184158-79-0P 184158-80-3P
     184158-81-4P 184158-82-5P 184158-83-6P
     184158-86-9P 184158-87-0P 184158-88-1P
     184377-51-3P
     RL: BAC (Biological activity or effector, except adverse); BSU (Biological
     study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
     BIOL (Biological study); PREP (Preparation); USES (Uses)
        (preparation of N-(oxobenzoxazol-6-yl)oxazolidinones and analogs as
        antibacterial agents)
RN
     184156-99-8 HCAPLUS
     Acetamide, N-[[3-(2,3-dihydro-3-methyl-2-oxo-6-benzothiazolyl)-2-oxo-5-
CN
     oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)
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RN 184157-00-4 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-1,3-dimethyl-2-oxo-1H-benzimidazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-01-5 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-3-methyl-1-(3-methylbutyl)-2-oxo-1H-benzimidazol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-02-6 HCAPLUS

CN Acetamide, N-[[3-[1-ethyl-2,3-dihydro-2-oxo-3-(phenylmethyl)-1H-benzimidazol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 184157-03-7 HCAPLUS

CN Acetamide, N-[[3-(3-ethyl-2,3-dihydro-2-oxo-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-06-0 HCAPLUS

CN Acetamide, N-[[3-(1-ethyl-2,3-dihydro-2-oxo-1H-benzimidazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-08-2 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-10-6 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2-oxo-5-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

184157-12-8 HCAPLUS RN

Acetamide, N-[[3-(2,3-dihydro-2-thioxo-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN . 184157-14-0 HCAPLUS

Acetamide, N-[[3-(2,3-dihydro-2-thioxo-5-benzoxazolyl).-2-oxo-5oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-17-3 HCAPLUS RN

Acetamide, N-[[3-(2,3-dihydro-2-thioxo-6-benzothiazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-19-5 HCAPLUS RN

Acetamide, N-[[3-(2-amino-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, CN (S) - (9CI)(CA INDEX NAME)

Absolute stereochemistry.

RN 184157-21-9 HCAPLUS

Acetamide, N-[[3-(2-amino-5-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, CN (S) - (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-23-1 HCAPLUS RN

Acetamide, N-[[3-(2-amino-6-benzoxazoly1)-2-oxo-5-oxazolidiny1]methy1]-,CN monohydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HC1

184157-26-4 HCAPLUS RN

Acetamide, N-[[3-(2-amino-5-benzoxazoly1)-2-oxo-5-oxazolidiny1]methy1]-, CN monohydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HC1

RN 184157-29-7 HCAPLUS

Acetamide, N-[[3-(3-acetyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-32-2 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-3-(methylsulfonyl)-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-35-5 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-[(phenylmethoxy)acetyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-38-8 HCAPLUS

CN Acetamide, N-[[3-[3-(cyclopropylcarbonyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-41-3 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-3-methyl-2-oxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-44-6 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-(phenylmethyl)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-45-7 HCAPLUS

CN 3(2H)-Benzoxazoleacetic acid, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-oxo-, ethyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-47-9 HCAPLUS

CN Acetamide, N-[[3-[3-(cyanomethyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-

5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-49-1 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-(2,2,2-trifluoroethyl)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-50-4 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-3-(hydroxymethyl)-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 184157-51-5 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-3-methyl-2-thioxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-52-6 HCAPLUS

CN Acetamide, N-[[3-(3-ethyl-2,3-dihydro-2-thioxo-6-benzoxazolyl)-2-oxo-5-

oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-53-7 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-3-propyl-2-thioxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-54-8 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-3-methyl-2-thioxo-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-55-9 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-3-(1-methylethyl)-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-56-0 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-3-(1-methylpropyl)-2-oxo-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (5S)-[partial]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-57-1 HCAPLUS RN

Acetamide, N-[[3-[3-(1-ethylpropyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-58-2 HCAPLUS RN

Acetamide, N-[[3-(3-cyclopropyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-59-3 HCAPLUS RN

Acetamide, N-[[3-(3-cyclobutyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 184157-60-6 HCAPLUS

CN Acetamide, N-[[3-(3-cyclopentyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-61-7 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-3-methyl-2-oxo-5-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-64-0 HCAPLUS

CN Propanamide, N-[[(5S)-3-(2,3-dihydro-3-methyl-2-oxo-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

184157-88-8 HCAPLUS

Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-(2-propenyl)-6-benzoxazolyl]-2-oxo-5oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-89-9 HCAPLUS

RN Acetamide, N-[[3-(3-ethyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-90-2 HCAPLUS RN

Acetamide, N-[[3-(2,3-dihydro-2-oxo-3-propyl-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-91-3 HCAPLUS RN

Acetamide, N-[[3-(3-butyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

184157-92-4 HCAPLUS RN

Acetamide, N-[[3-(2,3-dihydro-2-oxo-3-pentyl-6-benzoxazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-93-5 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-3-(2-methylpropyl)-2-oxo-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-94-6 HCAPLUS

Acetamide, N-[[3-[2,3-dihydro-3-(2-hydroxyethyl)-2-oxo-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184157-95-7 HCAPLUS RN

Acetamide, N-[[3-[3-[(dimethylamino)methyl]-2,3-dihydro-2-oxo-6-CN . benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-96-8 HCAPLUS
CN Acetamide, N-[[3-[3-[(dimethylamino)methyl]-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, hydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

●x HCl

RN 184157-97-9 HCAPLUS
CN Acetamide, N-[[3-(2,3-dihydro-2-oxo-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-98-0 HCAPLUS
CN Acetamide, N-[[3-[2,3-dihydro-3-(hydroxymethyl)-2-oxo-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184157-99-1 HCAPLUS

CN Acetamide, N-[[3-[3-(fluoromethyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-00-7 HCAPLUS

CN Acetamide, N-[[3-[3-(fluoromethyl)-2,3-dihydro-2-oxo-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-01-8 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-3-methyl-2-(2-propenylimino)-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

RN 184158-02-9 HCAPLUS

CN Cyclopropanecarboxamide, N-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-2(3H)-benzothiazolylidene]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

184158-03-0 HCAPLUS · RN

Propanamide, N-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-10-0x0-3-0x0-10-0CN 2(3H)-benzothiazolylidene]-2,2-dimethyl-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

184158-04-1 HCAPLUS RN

Acetamide, N-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-CN 2(3H)-benzothiazolylidene]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. . Double bond geometry unknown.

RN 184158-05-2 HCAPLUS

Acetamide, N-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-CN 2(3H)-benzothiazolylidene]-2,2,2-trichloro-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

184158-06-3 HCAPLUS RN

Carbamic acid, [6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-2(3H)-benzothiazolylidene]-, phenylmethyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

184158-07-4 HCAPLUS RN

Carbamic acid, [6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-CN 2(3H)-benzothiazolylidene]-, 4-nitrophenyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

184158-08-5 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-3-methyl-2-[[(methylamino)carbonyl]imino]-6-CN benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

RN 184158-09-6 HCAPLUS

Acetamide, N-[[3-[2-(cyanoimino)-2,3-dihydro-3-methyl-6-benzothiazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

184158-23-4 HCAPLUS RN

Propanamide, N-[[3-(3-ethyl-2,3-dihydro-2-oxo-6-benzothiazolyl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184158-29-0 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-3-(1-methylethyl)-2-oxo-6-benzothiazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184158-31-4 HCAPLUS RN

Propanamide, N-[[3-[2,3-dihydro-3-(1-methylethyl)-2-oxo-6-benzothiazolyl]-CN 2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-39-2 HCAPLUS CN Propanamide, N-[[3-[2,3-dihydro-3-(1-methylethyl)-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-54-1 HCAPLUS
CN Acetamide, N-[[3-[3-(2-chloro-1-methylethyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (5S)-[partial]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-56-3 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-[1-(phenylmethyl)-3-pyrrolidinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (5S)-[partial]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-58-5 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-[1-(phenylmethyl)-4-piperidinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-60-9 HCAPLUS

CN Propanamide, N-[[3-(2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-66-5 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-(4-pyridinylmethyl)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-67-6 HCAPLUS

CN Propanamide, N-[[3-(2,3-dihydro-3-methyl-2-oxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 184158-68-7 HCAPLUS

CN Propanamide, N-[[3-(3-ethyl-2,3-dihydro-2-oxo-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-69-8 HCAPLUS

CN Propanamide, N-[[3-(2,3-dihydro-2-oxo-3-propyl-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-70-1 HCAPLUS

CN Propanamide, N-[[3-[2,3-dihydro-3-(2-methylpropyl)-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 184158-71-2 HCAPLUS

CN 3(2H)-Benzoxazoleacetic acid, 2-oxo-6-[2-oxo-5-[[(1-oxopropyl)amino]methyl]-3-oxazolidinyl]-, phenylmethyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-72-3 HCAPLUS

CN Propanamide, N-[[3-[3-[2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)ethyl]-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-73-4 HCAPLUS

CN Propanamide, N-[[3-[3-[3-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)propyl]-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)-(9CI) (CA INDEX NAME)

RN 184158-78-9 HCAPLUS

CN Acetamide, N-[[3-[3-(2-azido-1-methylethyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (5S)-[partial]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-79-0 HCAPLUS

CN Propanamide, N-[[3-[3-(3-aminopropyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-80-3 HCAPLUS

CN Propanamide, N-[[3-[3-(2-aminoethyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

184158-81-4 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-2-oxo-3-(4-piperidinyl)-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

184158-82-5 HCAPLUS

RN 3(2H)-Benzoxazoleacetic acid, 6-[5-[(acetylamino)methyl]-2-oxo-3-CN oxazolidinyl]-2-oxo-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 184158-83-6 HCAPLUS

Acetamide, N-[[3-[3-(2-amino-1-methylethyl)-2,3-dihydro-2-oxo-6-methylethyl)]CN benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (5S)-[partial]- (9CI) INDEX NAME)

RN 184158-86-9 HCAPLUS
CN Propanamide, N-[[3-[3-(3-aminopropyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, hydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

●x HCl

RN 184158-87-0 HCAPLUS
CN Propanamide, N-[[3-[2,3-dihydro-2-oxo-3-(4-pyridinylmethyl)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, hydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

•x HCl

RN 184158-88-1 HCAPLUS
CN Propanamide, N-[[3-[2,3-dihydro-3-[(5-nitro-2-pyridinyl)methyl]-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, hydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

●x HCl

RN 184377-51-3 HCAPLUS

CN Acetamide, N-[[3-[3-(2-amino-1-methylethyl)-2,3-dihydro-2-oxo-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, hydrochloride, (5S)-[partial]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

●x HCl

IT 176490-94-1

RL: RCT (Reactant); RACT (Reactant or reagent)
 (preparation of N-(oxobenzoxazol-6-yl)oxazolidinones and analogs as
 antibacterial agents)

RN 176490-94-1 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[(phenylmethyl)thio]-6-benzothiazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

IT 176490-91-8P 176491-09-1P 184159-05-5P
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
 (Reactant or reagent)
 (preparation of N-(oxobenzoxazol-6-yl)oxazolidinones and analogs as
 antibacterial agents)
RN 176490-91-8 HCAPLUS
CN Acetamide, N-[[3-[2-(methylthio)-6-benzothiazolyl]-2-oxo-5 oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-09-1 HCAPLUS

CN Benzothiazolium, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-2-(methylthio)-, iodide, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

• I-

RN 184159-05-5 HCAPLUS
CN Acetamide, N-[[3-(2,3-dihydro-2-imino-3-methyl-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

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ACCESSION NUMBER: 1996:294873 HCAPLUS

DOCUMENT NUMBER: 124:343286

TITLE:

Preparation of (benzoxazolyl) - and

(benzothiazolyl)oxazolidinone antibiotics

INVENTOR(S):

Haebisch, Dieter; Riedl, Bernd; Ruppelt, Martin; Stolle, Andreas; Wild, Hanno; Endermann, Rainer;

Bremm, Klaus-Dieter; Kroll, Hein-Peter; Labischinski,

Harald; et al.

PATENT ASSIGNEE(S):

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EP 697412	В1	20000112		
R: AT, BE, CH,	DE, DK	, ES, FR,	GB, GR, IE, IT, LI, LU,	MC, NL, PT, SE
AT 188698	E	20000115	AT 1995-111477	
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IL 114784	A1	19991028	IL 1995-114784	19950731
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ZA 9506445	Α	19960322	ZA 1995-6445	19950802
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PRIORITY APPLN. INFO.:			DE 1994-4427475	A1 19940803
			DE 1995-19514313	A 19950418

OTHER SOURCE(S):

MARPAT 124:343286

GΙ

CN, CHO, CF3, NO2, (un)branched alkoxy, etc.; R1 = OSO2R3, tertiary-amino group; R3 = (un)substituted alkyl, (un)substituted Ph; R2 = H, CHO, CO2H, alkoxycarbonyl, etc.; * = chiral C], useful in pharmaceuticals, especially as antibiotics, are prepared Thus, benzothiazole II, m.p. 228°, was prepared by the m-chloroperbenzoic acid oxidation of the methylthio analog, and demonstrated a MIC against Staph. 133 (sic) of 1 $\mu g/mL$ and Staph. 48N (sic) of 2 mg/mL. 176490-91-8P 176490-92-9P 176490-93-0P ΙT 176490-94-1P 176490-95-2P 176490-96-3P 176490-97-4P 176490-98-5P 176490-99-6P 176491-00-2P 176491-01-3P 176491-02-4P 176491-03-5P 176491-04-6P 176491-05-7P 176491-06-8P 176491-07-9P 176491-08-0P 176491-09-1P 176491-13-7P 176491-14-8P 176491-15-9P 176491-24-0P 176491-25-1P 176491-26-2P 176491-27-3P 176491-28-4P 176491-29-5P 176491-30-8P 176491-31-9P 176491-32-0P 176491-33-1P 176491-34-2P 176491-35-3P 176491-36-4P 176491-37-5P 176491-38-6P 176491-39-7P 176491-40-0P · 176491-41-1P 176491-42-2P 176491-43-3P 176491-44-4P 176491-45-5P 176491-46-6P 176491-47-7P 176491-48-8P 176491-49-9P 176491-50-2P 176491-51-3P 176491-52-4P 176491-53-5P 176491-54-6P 176491-55-7P 176491-56-8P 176491-57-9P 176491-58-0P 176491-59-1P RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of (benzoxazolyl) - and (benzothiazolyl)oxazolidinone antibiotics) 176490-91-8 HCAPLUS RN Acetamide, N-[[3-[2-(methylthio)-6-benzothiazolyl]-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

The title compds. [I; A = O, S(O)a; a = O, 2; G, L, M = H, CO2H, halogen,

Absolute stereochemistry.

RN 176490-92-9 HCAPLUS
CN Acetamide, N-[[3-(2-methyl-6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-,
(S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176490-93-0 HCAPLUS CN Acetamide, N-[[3-(2-methyl-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176490-94-1 HCAPLUS
CN Acetamide, N-[[2-oxo-3-[2-[(phenylmethyl)thio]-6-benzothiazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176490-95-2 HCAPLUS
CN Acetamide, N-[[2-oxo-3-[2-(2-phenylethenyl)-6-benzothiazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry unknown.

RN 176490-96-3 HCAPLUS
CN Acetamide, N-[[2-oxo-3-(2-phenyl-5-benzoxazolyl)-5-oxazolidinyl]methyl]-,
(S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176490-97-4 HCAPLUS
CN Acetamide, N-[[3-(2-methyl-5-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176490-98-5 HCAPLUS
CN Benzothiazolium, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,3-dimethyl-, iodide, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

• I -

RN 176490-99-6 HCAPLUS
CN Acetamide, N-[[3-[2-(methylsulfinyl)-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

$$\begin{array}{c|c} & & & & \\ & & & \\ & & & \\ & & & \\ \text{AcNH-CH}_2 & & & \\ \end{array}$$

RN 176491-00-2 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[(phenylmethyl)sulfonyl]-6-benzothiazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 176491-01-3 HCAPLUS

CN Acetamide, N-[[3-[2-(cyclopropylamino)-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-02-4 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-(2-pyridinylthio)-6-benzothiazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-03-5 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[(phenylmethyl)sulfinyl]-6-benzothiazolyl]-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

AcNH-CH2

RN 176491-04-6 HCAPLUS

CN Acetamide, N-[[3-[2-(methylsulfonyl)-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 176491-05-7 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-(2-pyrimidinylthio)-6-benzothiazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-06-8 HCAPLUS

CN Acetamide, N-[[3-[2-(chloromethyl)-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-07-9 HCAPLUS

CN Acetamide, N-[[3-(2-formyl-6-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-08-0 HCAPLUS

CN Acetamide, N-[[3-[2-(hydroxymethyl)-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-09-1 HCAPLUS

CN Benzothiazolium, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-3-methyl-2-(methylthio)-, iodide, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

) I-

RN 176491-13-7 HCAPLUS

CN Acetamide, N-[[3-[2-[2-(4-chlorophenyl)ethenyl]-6-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry unknown.

RN 176491-14-8 HCAPLUS

CN Acetamide, N-[[3-[2-(3-butenyl)-5-benzothiazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

176491-15-9 HCAPLUS RN

Carbonothioic acid, S-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-CN benzothiazolyl] O-ethyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

176491-24-0 HCAPLUS RN

Acetamide, N-[[3-(6-benzoxazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)-CN (CA INDEX NAMÉ)

Absolute stereochemistry.

176491-25-1 HCAPLUS RN

Acetamide, N-[[3-[2-[4-(4-methoxyphenyl)-1-piperazinyl]-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

176491-26-2 HCAPLUS RN'

Acetamide, N-[[2-oxo-3-[2-[4-[3-(trifluoromethyl)phenyl]-1-piperazinyl]-6-CN benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 176491-27-3 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[4-[4-(trifluoromethyl)phenyl]-1-piperazinyl]-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-28-4 HCAPLUS

CN Acetamide, N-[[3-[2-[4-[2-(4-morpholinyl)-2-oxoethyl]-1-piperazinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-29-5 HCAPLUS

CN Acetamide, N-[[3-[2-[4-(2-methoxyethyl)-1-piperazinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-30-8 HCAPLUS

CN Acetamide, N-[[3-[2-(hexahydro-1H-1,4-diazepin-1-yl)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-31-9 HCAPLUS

CN Acetamide, N-[[3-[2-[[2-(4-methoxyphenyl)ethyl]amino]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-32-0 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[[[4-(trifluoromethyl)phenyl]methyl]amino]-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-33-1 HCAPLUS
CN 3-Piperidinecarboxamide, 1-[6-[5-[(acetylamino)methyl]-2-oxo-3oxazolidinyl]-2-benzoxazolyl]- (9CI) (CA INDEX NAME)

$$\begin{array}{c|c} & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & &$$

RN 176491-34-2 HCAPLUS
CN Pyridinium, 1-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-benzoxazolyl]-4-(dimethylamino)-, iodide, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

• I-

RN 176491-35-3 HCAPLUS
CN Acetamide, N-[[2-oxo-3-[2-(2-pyridinyloxy)-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-36-4 HCAPLUS
CN Acetamide, N-[[3-[2-[(2-hydroxyethyl)amino]-6-benzoxaźolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 176491-37-5 HCAPLUS
CN Acetamide, N-[[3-[2-[(3-hydroxypropyl)amino]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-38-6 HCAPLUS

CN Acetamide, N-[[3-[2-[[2-(2-hydroxyethoxy)ethyl]amino]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-39-7 HCAPLUS

CN Acetamide, N-[[3-[2-[3-(2-hydroxyethyl)-1-imidazolidinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-40-0 HCAPLUS

CN Acetamide, N-[[3-[2-[(2-hydroxypropyl)amino]-6-benzoxazolyl]]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 176491-41-1 HCAPLUS

CN Acetamide, N-[[3-[2-[(2-hydroxy-1,1-dimethylethyl)amino]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-42-2 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[4-(2-pyridinyl)-1-piperazinyl]-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-43-3 HCAPLUS

CN Acetamide, N-[[3-[2-[4-(1,3-benzodioxol-5-ylmethyl)-1-piperazinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 176491-44-4 HCAPLUS

CN Acetamide, N-[[3-[2-[4-[2-(4-morpholinyl)ethyl]-1-piperazinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-45-5 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-(4-pyrazinyl-1-piperazinyl)-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-46-6 HCAPLUS

CN Acetamide, N-[[3-[2-[4-(4-nitrophenyl)-1-piperazinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-47-7 HCAPLUS

CN Acetamide, N-[[3-[2-(cyclopropylmethoxy)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-48-8 HCAPLUS

CN Acetamide, N-[[3-[2-[4-[2-(dimethylamino)ethyl]-1-piperazinyl]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-49-9 HCAPLUS

CN Acetamide, N-[[3-[2-(1,4-dioxa-8-azaspiro[4.5]dec-8-yl)-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

176491-50-2 HCAPLUS RN

Acetamide, N-[[3-[2-(4-hydroxy-1-piperidinyl)-6-benzoxazolyl]-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

176491-51-3 HCAPLUS RN

Acetamide, N-[[2-oxo-3-[2-[[3-(1-piperidinyl)propyl]amino]-6-benzoxazolyl]-6-benzoxazolyl]CN 5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-52-4 HCAPLUS

Acetamide, N-[[3-[2-[(1,3-benzodioxol-5-ylmethyl)amino]-6-benzoxazolyl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

176491-53-5 HCAPLUS RN

Acetamide, N-[[2-oxo-3-[2-[(2-pyridinylmethyl)amino]-6-benzoxazolyl]-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 176491-54-6 HCAPLUS

CN Acetamide, N-[[3-[2-[[3-(dimethylamino)propyl]methylamino]-6-benzoxazolyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-55-7 HCAPLUS

CN 1-Piperazinecarboxylic acid, 4-[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-benzoxazolyl]-, 1,1-dimethylethyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-56-8 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-(1-piperazinyl)-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, monohydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

RN 176491-57-9 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-(4-oxo-1-piperidinyl)-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 176491-58-0 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[2-[(2-pyridinylmethyl)amino]-6-benzoxazolyl]-5-oxazolidinyl]methyl]-, monohydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

RN 176491-59-1 HCAPLUS

CN Propanamide, 2-[[6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-benzoxazolyl]amino]-N-methyl- (9CI) (CA INDEX NAME)

AcNH-CH2

L23 ANSWER 40 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1996:241534 HCAPLUS

DOCUMENT NUMBER:

124:289520

TITLE:

Preparation of 3-heteroaryl-2-oxazolidinones as

antibacterials.

INVENTOR(S):

Riedl, Bernd; Haebich, Dieter; Stolle, Andreas; Wild, Hanno; Endermann, Rainer; Bremm, Klaus Dieter; Kroll, Hein-Peter; Labischinski, Harald; Schaller, Klaus;

Werling, Hans-Otto

PATENT ASSIGNEE(S):

Bayer A.-G., Germany

SOURCE:

Eur. Pat. Appl., 62 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent German

LANGUAGE:

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 693491 R: AT, BE, C	A1 H, DE, DK	19960124 , ES, FR, G	EP 1995-110629 B, GR, IE, IT, LI,	19950707 LU, MC, NL, PT, SE
DE 4425613	A1	19960125	DE 1994-4425613	19940720
RO 115161	В1	19991130	RO 1995-1295	19950711
AU 9524988	A1	19960201	AU 1995-24988	19950713
AU 695661	В2	19980820		
CA 2154026	AA	19960121	CA 1995-2154026	19950717
US 5698574	Α	19971216	US 1995-503183	19950717
IL 114622	A1	19990714	IL 1995-114622	19950717
FI 9503476	Α	19960121	FI 1995-3476	19950718
JP 08053443	A2	19960227	JP 1995-203904	19950718
NO 9502866	A	19960122	NO 1995-2866	19950719
ZA 9506015	A	19960222	ZA 1995-6015	19950719
CN 1121919	A	19960508	CN 1995-108995	19950719
ни 74003	A2	19961028	ни 1995-2167	19950719
PRIORITY APPLN. INFO.:			DE 1994-4425613	A 19940720
OTHER SOURCE(S): GI		124:289520		

Title compds. [I; R1 = N3, OH, OR2, OSO3R3, NR4R5; R2 = acyl, protecting AB group; R3 = alkyl, (alkyl-substituted) Ph; R, R5 = H, alkyl, cycloalkyl, Ph, acyl; A = C-bonded (substituted) 5-membered heteroaryl which may be benzo- or naphtho-fused], were prepared Thus, 4-bromobenzo[b]thiophene-2carboxylic acid and Et3N in acetone at 0° were treated with iso-Bu chloroformate and then with NaN3 to give 76% 4-bromobenzo[b]thiophene-2carbonyl azide. The latter was refluxed 1 h with LiBr and Bu3PO in xylene using a water separator; (R)-glycidyl butyrate in xylene was added at reflux to give 30% (5R)-3-[4-bromobenzo[b]thiophenyl]-5butyryloxymethyloxazolidin-2-one. This was stirred with cesium carbonate in MeOH to give 61% title compound (II). I inhibited Staph. 133 with a min. inhibitory concentration of $0.25-16 \mu g/mL$.

175591-46-5P 175591-47-6P 175591-48-7P IT 175591-49-8P 175591-57-8P 175591-58-9P 175591-61-4P 175591-90-9P 175591-91-0P 175591-92-1P 175591-93-2P 175591-94-3P 175591-95-4P 175591-96-5P 175591-97-6P 175591-98-7P 175592-05-9P 175592-06-0P 175592-09-3P 175592-10-6P 175592-11-7P 175592-12-8P 175592-13-9P 175592-14-0P 175592-15-1P 175592-16-2P 175592-17-3P

175592-18-4P 175592-19-5P 175592-20-8P 175592-21-9P 175592-22-0P 175592-23-1P 175592-24-2P 175592-37-7P 175592-38-8P 175592-39-9P 175592-40-2P 175592-41-3P 175592-43-5P 175592-45-7P 175592-46-8P 175592-47-9P 175592-49-1P 175592-50-4P 175592-51-5P 175592-52-6P RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of 3-heteroaryl-2-oxazolidinones as antibacterials) 175591-46-5 HCAPLUS RN Acetamide, N-[[3-(4-bromobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl]-CN , (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175591-47-6 HCAPLUS
CN Acetamide, N-[[3-(5-bromobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-48-7 HCAPLUS
CN Acetamide, N-[[3-(6-bromobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175591-49-8 HCAPLUS

CN Acetamide, N-[[3-(7-bromobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (+).

RN 175591-57-8 HCAPLUS

CN Acetamide, N-[[3-(6-bromo-2-benzothiazolyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175591-58-9 HCAPLUS

CN Acetamide, N-[[3-(5-bromo-2-benzofuranyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175591-61-4 HCAPLUS

CN Acetamide, N-[[3-(2-benzofuranyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175591-90-9 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(4-phenylbenzo[b]thien-2-yl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175591-91-0 HCAPLUS

CN Acetamide, N-[[3-[4-(4-methylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-92-1 HCAPLUS

CN Acetamide, N-[[3-[4-(2-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-93-2 HCAPLUS

CN Acetamide, N-[[3-[4-(3-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-94-3 HCAPLUS

CN Acetamide, N-[[3-[4-(4-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-95-4 HCAPLUS

CN Acetamide, N-[[3-[4-(4-acetylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-96-5 HCAPLUS

CN Acetamide, N-[[3-[4-(4-fluorophenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-97-6 HCAPLUS

CN Acetamide, N-[[3-[4-(3-aminophenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175591-98-7 HCAPLUS

CN Acetamide, N-[[3-[4-(3-acetylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-05-9 HCAPLUS

CN Acetamide, N-[[3-[5-(3-acetylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-06-0 HCAPLUS

CN Acetamide, N-[[3-[5-(4-fluorophenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-09-3 HCAPLUS
CN Acetamide, N-[[3-[5-(4-acetylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-10-6 HCAPLUS

CN Acetamide, N-[[3-[5-(3-aminophenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, hydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

●x HCl

RN 175592-11-7 HCAPLUS

CN Acetamide, N-[[3-[5-(3-chloro-4-fluorophenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

175592-12-8 HCAPLUS RN

Acetamide, N-[[3-[5-[3,5-bis(trifluoromethyl)phenyl]benzo[b]thien-2-yl]-2-CN oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-13-9 HCAPLUS

Acetamide, N-[[3-[4-(5-formyl-3-thienyl)benzo[b]thien-2-yl]-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN175592-14-0 HCAPLUS

CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-15-1 HCAPLUS

CN Acetamide, N-[[3-[4-[3,5-bis(trifluoromethyl)phenyl]benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-16-2 HCAPLUS

CN Acetamide, N-[[3-[4-(4-methoxyphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-17-3 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(7-phenylbenzo[b]thien-2-yl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-18-4 HCAPLUS

CN Acetamide, N-[[3-[7-(3-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-19-5 HCAPLUS

CN Acetamide, N-[[3-[5-(4-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-20-8 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(5-phenylbenzo[b]thien-2-yl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 175592-21-9 HCAPLUS

CN Acetamide, N-[[3-[5-(3-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-22-0 HCAPLUS

CN Acetamide, N-[[3-[5-(5-acetyl-2-thienyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-23-1 HCAPLUS

CN Acetamide, N-[[3-[7-(4-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 175592-24-2 HCAPLUS CN Acetamide, N-[[3-[7-(4-fluorophenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-37-7 HCAPLUS
CN Acetamide, N-[[3-(5-fluorobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-38-8 HCAPLUS
CN Acetamide, N-[[3-(5-methylbenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).

RN 175592-39-9 HCAPLUS CN Acetamide, N-[[3-(5-cyanobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-40-2 HCAPLUS
CN Acetamide, N-[[3-(6-fluorobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-41-3 HCAPLUS
CN Acetamide, N-[[3-(6-chlorobenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-43-5 HCAPLUS
CN Benzoic acid, 4-[2-[5-[(acetylamino)methyl]-2-oxo-3-

oxazolidinyl]benzo[b]thien-5-yl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-45-7 HCAPLUS

CN Acetamide, N-[[3-[6-(4-formylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-46-8 HCAPLUS

CN Acetamide, N-[[3-[6-(4-acetylphenyl)benzo[b]thien-2-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-47-9 HCAPLUS

CN Acetamide, N-[[3-(6-[1,1'-biphenyl]-4-ylbenzo[b]thien-2-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 175592-49-1 HCAPLUS

CN Benzo[b]thiophene-4-carboxylic acid, 2-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, methyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-50-4 HCAPLUS

CN Benzo[b]thiophene-6-carboxylic acid, 2-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, methyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175592-51-5 HCAPLUS

CN Benzo[b]thiophene-5-carboxylic acid, 2-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, methyl ester, (S)- (9CI) (CA INDEX NAME)

175592-52-6 HCAPLUS RN

Acetamide, N-[(3-benzo[b]thien-2-yl-2-oxo-5-oxazolidinyl)methyl]-, (S)-benzo[b]thien-2-yl-2-oxo-5-oxazolidinyl)methyl]-, (S)-benzo[b]thien-2-yl-2-oxo-5-oxazolidinyl]methyl]-, (S)-benzo[b]thien-2-yl-2-oxo-5-oxazolidinyl]methyllandylmCN (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 41 OF 50

ACCESSION NUMBER:

1996:231381 HCAPLUS

DOCUMENT NUMBER:

124:261020

TITLE:

Preparation of benzofuranyl- and

benzothienyloxazolidinones as antibacterials.

INVENTOR(S):

Riedl, Bernd; Haebich, Dieter; Stolle, Andreas; Wild, HannoEndermann, Rainer; Endermann, Rainer; Bremm,

Klaus-Dieter; Kroll, Hein-Peter; Labischinski, Harald;

Schaller, Klaus; Werling, Hans-Otto

PATENT ASSIGNEE(S): SOURCE:

Bayer A.-G., Germany Ger. Offen., 34 pp.

CODEN: GWXXBX

DOCUMENT TYPE:

Patent

LANGUAGE:

German

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 4425609	A1	19960125	DE 1994-4425609	19940720
EP 694544	A1	19960131	EP 1995-110628	19950707
R: AT, BE, CH,	DE, DK	, ES, FR, G	B, GR, IE, IT, LI,	LU, MC, NL, PT, SE
CA 2154024	AA	19960121	CA 1995-2154024	19950717
. US 5684023	Α	19971104	US 1995-503116	19950717
JP 08041057	A2	19960213	JP 1995-205247	19950719
PRIORITY APPLN. INFO.:			DE 1994-4425609	A 19940720
OTHER SOURCE(S):	MARPAT	124:261020		
CI				

AB Title compds. [I; R1 = N3, OH, OR2, OSO2R3, NR4R5; R2 = acyl, protecting group; R3 = alkyl, (substituted) Ph; r4, R5 = H, cycloalkyl, Ph, protecting group, COR6; R6 = H, cycloalkyl, alkyl, alkoxy, Ph; A = O, S; D, E, G, L, M = H, CO2H, halo, cyano, mercapto, formyl, CF3, NO2, alkyl, alkoxy, alkoxycarbonyl, alkylthio, acyl, etc.], and their salts and S-oxides, were prepared Thus, 5-benzyloxycarbonylaminobenzo[b]thiophene (preparation given) and 1,10-phenanthroline in THF at -70° were treated with BuLi and then with (R)-glycidyl butyrate to give title compound (II). Selected I showed min. inhibitory concns. of 1-16 μg/mL against Staph.

175424-85-8P 175424-86-9P 175424-87-0P 175424-88-1P 175424-89-2P 175424-92-7P 175424-93-8P 175424-94-9P 175424-95-0P 175424-96-1P 175424-97-2P 175424-98-3P 175424-99-4P 175425-01-1P

175425-02-2P 175425-03-3P RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of benzofuranyl- and benzothienyloxazolidinones as antibacterials)

RN 175424-85-8 HCAPLUS

CN Acetamide, N-[(3-benzo[b]thien-5-yl-2-oxo-5-oxazolidinyl)methyl]-, (S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-86-9 HCAPLUS

CN Acetamide, N-[[3-(3-methylbenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-87-0 HCAPLUS

CN Benzo[b]thiophene-2-carboxylic acid, 5-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-88-1 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(2-phenylbenzo[b]thien-5-yl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-89-2 HCAPLUS

CN Propanamide, N-[[2-oxo-3-(2-phenylbenzo[b]thien-5-yl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-92-7 HCAPLUS

CN Acetamide, N-[[3-(2-methylbenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-93-8 HCAPLUS CN Acetamide, N-[[3-(2-bromobenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-94-9 HCAPLUS
CN Acetamide, N-[[3-(2-bromo-3-methylbenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-95-0 HCAPLUS

CN Acetamide, N-[[3-(3-bromo-2-methylbenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-96-1 HCAPLUS

CN Acetamide, N-[[3-(3-bromo-2-phenylbenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-97-2 HCAPLUS

CN Acetamide, N-[[3-[2-(4-formylphenyl)-3-methylbenzo[b]thien-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-98-3 HCAPLUS

CN Acetamide, N-[[3-[2-(4-formylphenyl)benzo[b]thien-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175424-99-4 HCAPLUS

CN Acetamide, N-[[3-[2-(4-acetylphenyl)benzo[b]thien-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175425-00-0 HCAPLUS

CN Acetamide, N-[[3-[2-(4-fluorophenyl)benzo[b]thien-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175425-01-1 HCAPLUS
CN Acetamide, N-[[2-oxo-3-[2-(3-pyridinyl)benzo[b]thien-5-yl]-5oxazolidinyl]methyl]-, monohydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HCl

RN 175425-02-2 HCAPLUS
CN Acetamide, N-[[3-(2-bromo-3-methyl-1,1-dioxidobenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175425-03-3 HCAPLUS
CN Acetamide, N-[[3-(1,1-dioxidobenzo[b]thien-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

L23 ANSWER 42 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER: 1996:228503 HCAPLUS

DOCUMENT NUMBER: 124:261021

TITLE: Six-membered nitrogen-containing heteroaryl

oxazolidinones useful as antibacterials

INVENTOR(S): Riedl, Bernd; Haebich, Dieter; Stolle, Andreas; Wild,

Hanno; Endermann, Rainer; Bremm, Klaus Dieter; Kroll, Hein-Peter; Labischinski, Harald; Schaller, Klaus;

Werling, Hans-Otto

PATENT ASSIGNEE(S): SOURCE:

Bayer A.-G., Germany Eur. Pat. Appl., 99 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent German

LANGUAGE:

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PA'	TENT NO.		KIND	DATE	APPLICATION NO.	DATE
EP	694543 R: AT.	BE, CH,	A1 DE, DK		EP 1995-110624 GB, GR, IE, IT, LI,	
DE	4425612	,	A1	19960404	DE 1994-4425612	19940720
	9524985		A1	19960201	AU 1995-24985	19950713
	699940		В2	19981217		
	115262		В1	19991230	RO 1995-1312	19950714
	2154025		AA	19960121	CA 1995-2154025	19950717
JP	08041056		A2	19960213	` JP 1995-201799	19950717
_	5627181		Α	19970506	US 1995-503369	[.] 19950717
	114626		. A1	19990817	IL 1995-114626	19950717
	9503477		A	19960121	FI 1995-3477	19950718
NO	9502865	•	Α	19960122	NO 1995-2865 ·	19950719
	9506018		Α	19960313	ZA 1995-6018	19950719
HU	75035		A2	19970328	ни 1995-2173	19950719
	1119647		A	19960403	CN 1995-107584	19950720
US	5843967		Α	19981201	US 1996-749581	19961115
	Y APPLN.	INFO.:			DE 1994-4425612	A 19940720
	 ,-				US 1995-503369	A1 19950717
OTHER S	OURCE(S):		MARPAT	124:2610	21	•

GI

The title compds. I are prepared [in which R1 = N3, OH, OR2, OSO2R3, NR4R5; AB

R2 = acyl, protecting group; R3 = alkyl, Ph, alkylphenyl; R4, R5 =
 cycloalkyl, H, Ph, alkyl, protecting group, COR6; R6 = cycloalkyl, alkyl,
 Ph, H; D = [all optionally substituted] C-bound, 6-membered, aromatic,
 N-containing heterocyclyl; or 6-membered-ring-containing bi- or tricyclic
aromatic

N-containing heterocyclyl; or β -carbolin-3-yl; or indolizinyl]. For example, cyclization of 5-bromo-2-isocyanatopyridine-HCl with (R)-glycidyl butyrate in the presence of Bu3P:O in xylene gave 26% title compound II. This was subjected to a sequence of ester methanolysis (69%), conversion of the resulting alc. to a mesylate (95%), then to an azide (95%), and then to an amine (85%), N-acetylation of the amine (98%), and PdO-catalyzed coupling with 4-MeC6H4B(OH)2 (60%), to give title compound III [R = Me]. The analogously prepared compound III [R = H] had MIC of 0.5-1 μ /mL against 4 strains of Staphylococcus.

IT 175391-91-0P 175391-92-1P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)

(preparation of heteroaryloxazolidinones as antibacterials)

RN 175391-91-0 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(2-quinolinyl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175391-92-1 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(6-quinolinyl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

IT 175391-93-2P 175391-95-4P 175391-96-5P
175391-97-6P 175391-98-7P 175392-15-1P
175392-16-2P 175392-17-3P 175392-18-4P
175392-19-5P 175392-20-8P 175392-21-9P
175392-22-0P 175392-24-2P 175392-25-3P
175392-26-4P 175392-27-5P 175392-63-9P
175392-64-0P 175392-65-1P 175392-66-2P
175392-67-3P 175392-90-2P 175392-91-3P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);

Absolute stereochemistry.

RN 175391-95-4 HCAPLUS CN Acetamide, N-[[3-(3-methyl-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175391-96-5 HCAPLUS
CN Acetamide, N-[[3-(3-bromo-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-,
(S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175391-97-6 HCAPLUS
CN Acetamide, N-[[3-(6-bromo-2-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-,
(S)- (9CI) (CA INDEX NAME)

RN 175391-98-7 HCAPLUS

CN Acetamide, N-[[3-(8-bromo-2-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-15-1 HCAPLUS

CN Acetamide, N-[[3-[6-(5-methoxy-2-thienyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-16-2 HCAPLUS

CN Acetamide, N-[[3-[6-(3-formylphenyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-17-3 HCAPLUS

CN Acetamide, N-[[3-[6-(4-methylphenyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-18-4 HCAPLUS

CN Acetamide, N-[[3-[8-(3-formylphenyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-19-5 HCAPLUS

CN Acetamide, N-[[3-[8-[3,5-bis(trifluoromethyl)phenyl]-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-20-8 HCAPLUS

CN Acetamide, N-[[3-[8-(5-acetyl-2-thienyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-21-9 HCAPLUS

CN Acetamide, N-[[3-[8-(4-methylphenyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-22-0 HCAPLUS

CN Acetamide, N-[[3-[6-(4-acetylphenyl)-2-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-24-2 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(6-quinolinyl)-5-oxazolidinyl]methyl]-, monohydrochloride, (S)- (9CI) (CA INDEX NAME)

● HCl

RN 175392-25-3 HCAPLUS
CN Quinolinium, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1-methyl-, iodide, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

• I-

RN 175392-26-4 HCAPLUS
CN Quinolinium, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-1-ethyl-, iodide, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

● т-

RN 175392-27-5 HCAPLUS
CN Acetamide, N-[[3-(1-oxido-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-,
(S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-63-9 HCAPLUS

CN Acetamide, N-[[3-[3-(4-fluorophenyl)-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-64-0 HCAPLUS

CN Acetamide, N-[[3-[3-(4-formylphenyl)-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-65-1 HCAPLUS

CN Acetamide, N-[[3-[3-(4-acetylphenyl)-6-quinolinyl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-66-2 HCAPLUS CN Acetamide, N-[[2-oxo-3-(3-phenyl-6-quinolinyl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-67-3 HCAPLUS
CN Acetamide, N-[[2-oxo-3-[3-(3-pyridinyl)-6-quinolinyl]-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 175392-90-2 HCAPLUS
CN Acetamide, N-[[3-(3-bromo-6-quinolinyl)-2-oxo-5-oxazolidinyl]methyl]-,
monohydrochloride, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

HC1

HC1

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 43 OF 50

ACCESSION NUMBER:

1995:858608 HCAPLUS

DOCUMENT NUMBER:

123:256757

TITLE:

Preparation of indolo[2,1-b]quinazoline-6,12-dione

tuberculostatics

INVENTOR(S):

Baker, William R.; Mitscher, Lester A.

PATENT ASSIGNEE(S):

Pathogenesis Corp., USA

SOURCE:

PCT Int. Appl., 96 pp.

CODEN: PIXXD2

DOCUMENT TYPE:

Patent

LANGUAGE:

GΙ

English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.			KIND DATE			APPLICATION NO.					DATE						
 WO	WO 9513807		A1 19950526		WO 1994-US13259					19941117							
	W:		AT,	AU,	BB,	BG,	BR,	BY,	CA,	CH	, CN,	CZ,	DE,	DK,	EE,	ES,	FI,
		GB,	GE,	HU,	JP,	KE,	KG,	KP,	KR,	ΚZ	, LK,	LR,	LT,	LU,	LV,	MD,	MG,
		MN,	MW,	NL,	NO,	NZ,	PL,	PT,	RO,	RU	, SD,	SE,	SI,	SK,	ТJ,	TT,	UA,
		US,															
	RW:	KE,	MW,	SD,	SZ,	AT,	BE,	CH,	DE,	DK	, ES,	FR,	GB,	GR,	ΙE,	ΙT,	LU,
	•	MC,	NL,	PT,	SE,	BF,	BJ,	CF,	CG,	CI	, CM,	GA,	GN,	ML,	MR,	ΝE,	SN,
		TD,	TG														
US	5441	955			Α		1995	0815		US :	1993-	1547	84		_	9931	
AU	9512	100			A1		1995	0606		AU	1995-	1210	0			9941	
PRIORITY	APP	LN.	INFO	. :						US	1993-	1547	84			9931	
										WO	1994-	US13	259	1	₩ 1	9941	117
OTHER SC	DURCE	(S):			MAR	PAT	123:	2567	57								•

Ι

The title compds. [I; A-H = C, N; or A and B or C and D can be taken together to be N or S; R1-R4, R8, R10 = H, halogen, alkyl, cycloalkyl, (un)substituted heterocyclyl, (un)substituted amino, NO2, CN, CHO, etc.; R7, R9 = H, halogen, (un)substituted alkyl, cycloalkyl, (un)substituted heterocyclyl] useful for the treatment of multidrug-resistant Mycobacterium tuberculosis and M. leprae, are prepared Thus, 5-fluoroisatin was added to a solution of Me3COK and N-methylpyrrolidone, producing 8-fluoroindolo[2,1-b]quinazoline-6,12-dione, II, m.p. 273-276°, which demonstrated a MIC against multiple drug-resistant M. tuberculosis (10038) of <1 $\mu g/mL$, vs. 10 $\mu g/mL$ for tryptanthrin.

IT 169038-12-4P 169038-13-5P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of indolo[2,1-b]quinazoline-6,12-dione tuberculostatics)

RN 169038-12-4 HCAPLUS

CN Acetamide, N-[[3-(8-fluoro-6,12-dihydro-6,12-dioxoindolo[2,1-b]quinazolin-3-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 169038-13-5 HCAPLUS

CN Acetamide, N-[[3-(8-fluoro-6,12-dihydro-6,12-dioxoindolo[2,1-b]quinazolin-2-y1)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

L23 ANSWER 44 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1993:649939 HCAPLUS

DOCUMENT NUMBER:

119:249939

TITLE:

Antidepressants containing naphthyloxazolidones Nakai, Hideo; Yamada, Koichiro; Nomura, Sumihiro;

Matsumoto, Mamoru; Iwata, Hiroshi

PATENT ASSIGNEE(S):

SOURCE:

Tanabe Seiyaku Co, Japan

Jpn. Kokai Tokkyo Koho, 17 pp.

CODEN: JKXXAF

DOCUMENT TYPE:

INVENTOR(S):

LANGUAGE:

Patent Japanese

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 05155772	A2	19930622	JP 1992-151103	.19920423
JP 2551298	B2	19961106		
PRIORITY APPLN. INFO.:			JP 1991-191353	A1 19910425
OTHER SOURCE(S):	MARPAT	119:249939		
GI				

$$R^{1}$$
 N O R^{2}

Antidepressants contain naphthyloxazolidones I [R1 = H, OH, NO2, NH2, SO3H, aminosulfonyl, lower alkenyloxy, lower alkynyloxy, mono- or di-lower alkylaminocarbonyloxy, lower alkanoyloxy, lower alkoxy (which may be substituted with aryl, cycloalkyl, O-containing monoheterocyclyl, OH, lower alkoxy, cyano, di-lower alkylamino, aminocarbonyl, lower alkoxycarbonyl, lower alkanoyloxy, lower alkylthio, lower alkylsufinyl, lower alkylsulfonyl); R2 = OH, lower alkoxy, lower alkylsufonyloxy, N3, (lower alkyl- or alkanoyl-substituted) NH2] or their pharmacol. acceptable salts as active ingredients. Condensation of 20.0 g 2-naphthylamine with 16.7 g ethoxycarbonyl chloride in CH2Cl2 at room temperature overnight gave 25.26 g N-ethoxycarbonyl-2-naphthylamine, which (3.82 g) was refluxed with 3.13 g 2-(methoxymethyl)oxirane and Et3N for 3.5 h to afford 3.17 g 3-(2-naphthyl)-5-methoxymethyl-2-oxazolidone. The product at 10-7 M inhibited 78.4% MAO-A.

IT 135205-29-7P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of, as antidepressant)

135205-29-7 HCAPLUS

RN Acetamide, N-[[3-(2-naphthalenyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) CN INDEX NAME)

HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 45 OF 50

ACCESSION NUMBER:

1992:151623 HCAPLUS

DOCUMENT NUMBER:

116:151623

TITLE:

Antibacterials. Synthesis and structure-activity

studies of 3-aryl-2-oxooxazolidines. 4. Multiply-substituted aryl derivatives

AUTHOR(S):

Park, Chung Ho; Brittelli, David R.; Wang, C. L. J.; Marsh, Frank D.; Gregory, Walter A.; Wuonola, Mark A.; McRipley, Ronald J.; Eberly, Virginia S.; Slee, Andrew

M.; Forbes, Martin

CORPORATE SOURCE:

Drug Discovery Res., Chem. Sci., Exp. Stn., Du Pont Merck Pharm. Co., Wilmington, DE, 19880-0353, USA Journal of Medicinal Chemistry (1992), 35(6), 1156-65

SOURCE:

CODEN: JMCMAR; ISSN: 0022-2623

DOCUMENT TYPE:

Journal

LANGUAGE:

English

GI

AB The synthesis and structure-activity relationship (SAR) studies of the effect of different polysubstitution patterns in the aromatic ring of 5-(acetamidomethyl)oxazolidinones, e.g., (I Ar = substituted Ph, 5-indolyl, β -naphthyl) on antibacterial activity are presented. Compds. I were prepared by the six-step synthesis described previously via electrophilic aromatic substitution reactions of 3-substituted compds., and functional-group interchange reactions of 3,4-disubstituted compds. Antibacterial evaluation of compds. I against Staphylococcus aureus and Enterococcus faecalis gave the following results. The 2,4- and 2,5-disubstituted derivs. have weak or no antibacterial activity. Antibacterial activities of 3,4-disubstituted compds. are comparable to those of the 4-monosubstituted analogs for small 3-substituents (smaller than Br), but decline rapidly for larger 3-substituents. 3,4-Annulated derivs. are comparable in activity to their open-chain analogs. 3,5-Disubstituted and 3,4,5- and 2,4,6-trisubstituted derivs. are devoid of antibacterial activity.

Absolute stereochemistry.

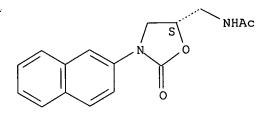
RN 120912-43-8 HCAPLUS
CN Acetamide, N-[[3-(2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 139071-73-1 HCAPLUS CN Acetamide, N-[[3-(1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 139071-74-2 HCAPLUS CN Acetamide, N-[[(5S)-3-(2-naphthalenyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)



HCAPLUS COPYRIGHT 2006 ACS on STN L23 ANSWER 46 OF 50

ACCESSION NUMBER:

1991:471581 HCAPLUS

DOCUMENT NUMBER:

115:71581

TITLE:

Preparation of naphthyloxazolidone derivatives as

selective monoamine oxidase A inhibitors

INVENTOR(S):

Nakai, Hideo; Yamada, Koichiro; Nomura, Sumihiro;

Matsumoto, Mamoru; Iwata, Hiroshi Tanabe Seiyaku Co., Ltd., Japan

PATENT ASSIGNEE(S):

Eur. Pat. Appl., 30 pp.

SOURCE:

CODEN: EPXXDW

DOCUMENT TYPE:

Patent

LANGUAGE:

English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 425209	A2	19910502	EP 1990-311524	19901019
EP 425209	A3	19920219		
EP 425209	B1	19950510	•	
R: AT, BE, CH,	DE, DK	, ES, FR, GB	B, GR, IT, LI, LU, NL,	
US 5182296	Α	19930126	US 1990-599564	19901018
AT 122344	Ε.	19950515	AT 1990-311524	19901019
ES 2074544	Т3	19950916	ES 1990-311524	19901019
CA 2028440	AA	19910427	CA 1990-2028440	19901024
CA 2028440	С	19961112		
JP 03218367	A2	19910925	JP 1990-289910	19901025
KR 128288	В1	19980402	KR 1990-17252	19901026
US 5332754	Α	19940726	US 1992-961873.	19921015
PRIORITY APPLN. INFO.:			JP 1989-279305	A 19891026
			US 1990-599564	A1 19901018
OMURD COURCE (C)	ייי א כו כו א א	115.71501	*	

OTHER SOURCE(S):

MARPAT 115:71581

GI

$$R^1$$
 N O R^2 T

Title compds. [I; R1 = H, OH, NO2, amino, sulfo, aminosulfonyl, AΒ alkenyloxy, alkynyloxy, alkylaminocarbonyloxy, alkanoyloxy, (substituted) . alkoxy; R2 = OH, alkoxy, alkylsulfonyloxy, triazo, amino], were prepared Thus, EtO2CC1, 2-naphthylamine, NaHCO3, and CH2C12 were stirred overnight to give N-ethoxycarbonyl-2-naphthylamine. The latter was refluxed with 2-methoxymethyloxirane and Et3N to give 5-methoxymethyl-3-(2-naphthyl)-2oxazolidene. I at 10-7 M gave 76.6-98.8% inhibition of MAO-A (monoamine oxidase A) from rat cerebral mitochondria; they were selective for MAO-A over MAO-B.

IT 135205-29-7P

RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of, as monoamine oxidase A inhibitor)

RN 135205-29-7 HCAPLUS

CN Acetamide, N-[[3-(2-naphthalenyl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

L23 ANSWER 47 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1990:572004 HCAPLUS

DOCUMENT NUMBER:

113:172004

TITLE:

Preparation of 5β -amidomethyloxazolidin-2-ones as

antibacterial agents Brickner, Steven L.

INVENTOR(S):
PATENT ASSIGNEE(S):

Upjohn Co., USA

SOURCE:

Eur. Pat. Appl., 75 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent

LANGUAGE:

English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATE	NT NO.			KINE	D DATE	APPLICATION NO.	DATE
EP 3	 59418 59418 R: ES,			A1 B1	19900321 19941012	EP 1989-308506	19890822
WO 90	002744			A1	19900322	WO 1989-US3548	19890822
Ţ	W: AU,	DK,	FI,	HU,	JP, KR, NO,	SU, US	•
AU 8	941957			A1	19900402 19911205	LU, NL, SE AU 1989-41957	19890822
						EP 1989-909990	19890822
		DF	CU	DE	ED CR TT	LT LU NL SE	
JP 0	4500665			T2	19920206	JP 1989-509255	19890822
				B2	19990308		
						EP 1994-102762	19890822
	09905				20010606		
		BE,	CH,	DE,	20010901	GR, IT, LI, LU, NL, SE ES 1994-102762	19890822
	157934			13 11	19950404	CA 1989-609594	
	38262			R1	19980515		
	164510				19921117		
	104515				19910313		
					19930126		19920423
	225565				19930706		
JP 1	1080139						19980707

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B2
                                 20010716
     JP 3188418
                                                                       20010831
                                              GR 2001-401350
                                  20011130
     GR 3036491
                           Т3
                                                                       20020606
                                              LV 2002-104
                                  20021220
                           В
     LV 12888
                                              DK 2004-1246
                                                                       20040818
                                  20040818
                           Α5
     DK 2004001246
                                  20050801
                           В1
     DK 175940
                                              US 1988-244988
                                                                      19880915
PRIORITY APPLN. INFO .:
                                              US 1988-253850
                                                                       19881005
                                                                   Α
                                              US 1989-324942
                                                                       19890317
                                              EP 1989-308506
                                                                   A3 19890822
                                              JP 1989-509255
                                                                   A3 19890822
                                              WO 1989-US3548
                                                                       19890822
                                                                   Α
                                              US 1991-655419
                                                                   A3
                                                                      19910220
                                              DK 1991-455
                                                                       19910313
                          MARPAT 113:172004
OTHER SOURCE(S):
     For diagram(s), see printed CA Issue.
GI
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The title compds. [I; W1, W2 = H, alkyl, etc., or W1W2 = NR5CH:CR7, NR5CH2CH2; R2-R5, R7 = H, alkyl, acyl, etc.; R6 = amido, e.g. NHAc], useful as antibacterials against gram-pos. and anaerobic infections, and against Mycobacterium avium in AIDS patients (no data), were prepared 3-(1-Acetyl-1H-indolin-5-yl)-5-(azidomethyl)-2-oxazolidinone (preparation given) was hydrogenated over Pd/C and the resulting amine acetylated with Ac2O to give 3-(1-acetyl-1H-indolin-5-yl)-5-(acetamidomethyl)-2-oxazolidinone.

IT 129487-98-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(preparation and reaction of, in preparation of antibacterials)

RN 129487-98-5 HCAPLUS

CN 1H-Indole-1-carboxylic acid, 5-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,3-dihydro-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

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ΙT
     120912-44-9P 129487-31-6P 129487-32-7P
     129487-33-8P 129487-34-9P 129487-35-0P
     129487-36-1P 129487-37-2P 129487-38-3P
     129487-39-4P 129487-40-7P 129487-41-8P
     129487-42-9P 129487-43-0P 129487-44-1P
     129487-45-2P 129487-46-3P 129487-47-4P 129487-48-5P 129487-49-6P 129487-52-1P
     129487-53-2P 129487-54-3P 129487-55-4P
     129487-56-5P 129487-57-6P 129487-58-7P
     129487-59-8P 129487-60-1P 129487-61-2P
     129487-62-3P 129487-63-4P 129487-64-5P
     129487-65-6P 129487-66-7P 129487-67-8P
     129487-73-6P 129487-74-7P 129487-75-8P
     129487-76-9P 129487-77-0P 129487-78-1P
     129487-79-2P 129487-80-5P 129487-81-6P
     129487-82-7P 129487-83-8P 129487-84-9P
     129487-85-0P 129487-86-1P 129491-47-0P
     129502-53-0P 129567-77-7P 129567-78-8P
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129567-79-9P 129567-80-2P 129567-81-3P 129567-82-4P 129567-83-5P 129567-84-6P 129567-85-7P 129567-86-8P 129567-87-9P 129567-88-0P 129567-89-1P 129567-90-4P 129567-91-5P 129567-95-9P 129567-96-0P 129567-98-2P 129567-99-3P 129568-00-9P 129568-01-0P 129568-02-1P 129568-71-4P 129568-72-5P 129645-53-0P 129954-20-7P 129954-21-8P RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of, as antibacterial) 120912-44-9 HCAPLUS RN Acetamide, N-[[3-(2,3-dihydro-1-hydroxy-1H-inden-5-yl)-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129487-31-6 HCAPLUS
CN Acetamide, N-[[3-(1-acetyl-2,3-dihydro-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-32-7 HCAPLUS
CN Acetamide, N-[[3-(2,3-dihydro-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl], (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-33-8 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(2-methylpropyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-34-9 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(1-oxopropyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-35-0 HCAPLUS

CN Acetamide, N-[[3-[1-(cyclopentylcarbonyl)-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-36-1 · HCAPLUS

CN Acetamide, N-[[3-(1-formyl-2,3-dihydro-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-37-2 HCAPLUS

CN Acetamide, N-[[3-[1-(dichloroacetyl)-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-38-3 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(phenylacetyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-39-4 HCAPLUS

CN Acetamide, N-[[3-[1-[(acetyloxy)acetyl]-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

129487-40-7 HCAPLUS · RN

Acetamide, N-[[3-[2,3-dihydro-1-(2-thienylcarbonyl)-1H-indol-5-yl]-2-oxo-5-indol-5-yl]CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

129487-41-8 HCAPLUS RN

Acetamide, N-[[3-(1-benzoyl-2,3-dihydro-1H-indol-5-yl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

129487-42-9 HCAPLUS RN

Acetamide, N-[[3-(2,3-dihydro-1-methyl-1H-indol-5-yl)-2-oxo-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 129487-43-0 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(hydroxyacetyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-44-1 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-[(phenylmethoxy)acetyl]-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-45-2 HCAPLUS

CN Acetamide, N-[[3-[1-(4-chlorobenzoyl)-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 129487-46-3 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(2-propenyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-47-4 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-1-propyl-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-48-5 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(methoxyacetyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 129487-49-6 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(1-oxohexyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-52-1 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2-methyl-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-53-2 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2-methyl-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

RN 129487-54-3 HCAPLUS

CN Acetamide, N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-55-4 HCAPLUS

CN Acetamide, N-[[3-(1',3'-dihydro-1'-oxospiro[cyclopropane-1,2'-[2H]inden]-5'-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-56-5 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(5,6,7,8-tetrahydro-6-methyl-5-oxo-2-naphthalenyl)-5-oxazolidinyl]methyl]-, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-57-6 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(5,6,7,8-tetrahydro-6-methyl-5-oxo-2-naphthalenyl)-5-oxazolidinyl]methyl]-, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-58-7 HCAPLUS

CN Acetamide, N-[[3-(3',4'-dihydro-1'-oxospiro[cyclopropane-1,2'(1'H)-naphthalen]-6'-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-59-8 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-(hydroxymethyl)-1-oxo-1H-inden-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-60-1 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-2-(hydroxymethyl)-1-oxo-1H-inden-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

RN 129487-61-2 HCAPLUS
CN Acetamide, N-[[2-oxo-3-[5,6,7,8-tetrahydro-6-(hydroxymethyl)-5-oxo-2-naphthalenyl]-5-oxazolidinyl]methyl]-, [S-(R*,S*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-62-3 HCAPLUS

CN Acetamide, N-[[2-oxo-3-[5,6,7,8-tetrahydro-6-(hydroxymethyl)-5-oxo-2-naphthalenyl]-5-oxazolidinyl]methyl]-, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-63-4 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2,2-dimethyl-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-64-5 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(5,6,7,8-tetrahydro-6,6-dimethyl-5-oxo-2-naphthalenyl)-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-65-6 HCAPLUS

CN 1H-Indole-1-carboxylic acid, 6-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2,3-dihydro-, 1,1-dimethylethyl ester (9CI) (CA INDEX NAME)

AcNH-CH2

RN 129487-66-7 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-1H-indol-6-yl)-2-oxo-5-oxazolidinyl]methyl](9CI) (CA INDEX NAME)

AcNH-CH2

RN 129487-67-8 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(2-propenyl)-1H-indol-6-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

AcNH-CH2

RN 129487-73-6 HCAPLUS

Acetamide, N-[[3-[2,3-dihydro-1-(hydroxyimino)-1H-inden-5-yl]-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129487-74-7 HCAPLUS

Acetamide, N-[[2-oxo-3-(5,6,7,8-tetrahydro-2-naphthalenyl)-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

129487-75-8 HCAPLUS RN

Acetamide, N-[[2-oxo-3-(5,6,7,8-tetrahydro-5-oxo-2-naphthalenyl)-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129487-76-9 HCAPLUS

Acetamide, N-[[3-(1H-indazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI)CN (CA INDEX NAME)

129487-77-0 HCAPLUS RN

Acetamide, N-[[3-(1-ethyl-1H-indazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-CN (CA INDEX NAME)

129487-78-1 HCAPLUS RNCN

Acetamide, N-[[3-(1-acetyl-1H-indazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-(CA INDEX NAME)

RN 129487-79-2 HCAPLUS

Acetamide, N-[[2-oxo-3-(1-propyl-1H-indazol-5-yl)-5-oxazolidinyl]methyl]-CN (CA INDEX NAME)

$$\begin{array}{c|c} & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ & & \\ &$$

129487-80-5 HCAPLUS RN

Acetamide, N-[[3-(1-ethyl-2-methyl-1H-benzimidazol-5-yl)-2-oxo-5-CNoxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

129487-81-6 HCAPLUS RN

Acetamide, N-[[2-oxo-3-(1-propyl-1H-benzimidazol-5-yl)-5-CN oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-82-7 HCAPLUS

CN 1H-Benzimidazole-1-carboxylic acid, 5-[5-[(acetylamino)methyl]-2-oxo-3-oxazolidinyl]-2-methyl-, 1,1-dimethylethyl ester, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-83-8 HCAPLUS

CN Acetamide, N-[[3-(2-methyl-1H-benzimidazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-84-9 HCAPLUS

CN Acetamide, N-[[3-(1-acetyl-2-methyl-1H-benzimidazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-85-0 HCAPLUS
CN Acetamide, N-[[3-(1-formyl-1H-benzimidazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129487-86-1 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(methylsulfonyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129491-47-0 HCAPLUS

CN Acetamide, N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, [S-(R*,R*)]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129502-53-0 HCAPLUS

CN Acetamide, N-[[3-[1-(chloroacetyl)-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

129567-77-7 HCAPLUS RN

Acetamide, N-[[3-(1-acetyl-2,3-dihydro-1H-indol-5-yl)-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-78-8 HCAPLUS

Acetamide, N-[[3-(2,3-dihydro-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-CN (CA INDEX NAME)

AcNH-CH2

RN 129567-79-9 HCAPLUS

Acetamide, N-[[3-[2,3-dihydro-1-(2-methylpropyl)-lH-indol-5-yl]-2-oxo-5-indol-5-yl]CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

129567-80-2 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-1-(2-propenyl)-1H-indol-5-yl]-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Acnh-CH₂
$$CH_2$$
-CH=CH₂

RN 129567-81-3 HCAPLUS CN Acetamide, N-[[3-(2,3-dihydro-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

AcNH-CH2

RN 129567-82-4 HCAPLUS
CN Acetamide, N-[[3-(2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-83-5 HCAPLUS CN Acetamide, N-[[3-(1H-indazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)-(9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129567-84-6 HCAPLUS
CN Acetamide, N-[[2-oxo-3-(1-propyl-1H-indazol-5-yl)-5-oxazolidinyl]methyl]-,
(S)- (9CI) (CA INDEX NAME)

RN 129567-85-7 HCAPLUS

CN Acetamide, N-[[3-(1-benzoy1-2,3-dihydro-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

Acnh-CH₂
$$C-Ph$$

RN 129567-86-8 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-1-methyl-1H-indol-5-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-87-9 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-2,2-dimethyl-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-88-0 HCAPLUS

CN Acetamide, N-[[3-(1',3'-dihydro-1'-oxospiro[cyclopropane-1,2'-[2H]inden]-5'-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-89-1 HCAPLUS

CN Acetamide, N-[[2-oxo-3-(5,6,7,8-tetrahydro-6,6-dimethyl-5-oxo-2-naphthalenyl)-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-90-4 HCAPLUS

CN Acetamide, N-[[3-[1-[(acetyloxy)acetyl]-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-91-5 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(2-thienylcarbonyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-95-9 HCAPLUS

CN Acetamide, N-[[3-(1-ethyl-2-methyl-1H-benzimidazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-96-0 HCAPLUS

CN Acetamide, N-[[3-(1-ethyl-1H-indazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 129567-98-2 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(hydroxyacetyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129567-99-3 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-[(phenylmethoxy)acetyl]-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129568-00-9 HCAPLUS

CN Acetamide, N-[[3-[1-(4-chlorobenzoyl)-2,3-dihydro-1H-indol-5-yl]-2-oxo-5-

oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129568-01-0 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(methoxyacetyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129568-02-1 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(1-oxohexyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129568-71-4 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-(methylsulfonyl)-1H-indol-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

129568-72-5 HCAPLUS RN

Acetamide, N-[[3-(1-acetyl-1H-indazol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, CN (S) - (9CI) (CA INDEX NAME)

Absolute stereochemistry.

129645-53-0 HCAPLUS RN

Acetamide, N-[[3-[2,3-dihydro-1-propyl-1H-indol-5-yl]-2-oxo-5-CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

129954-20-7 HCAPLUS RN

Acetamide, N-[[3-(2,3-dihydro-2-methyl-1-oxo-1H-inden-5-yl)-2-oxo-5-methyl-1-oxo-1H-inden-5-wethyl-1-oxo-1H-indeCN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 129954-21-8 HCAPLUS

Acetamide, N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-kernel N-[[3-(2-ethyl-2,3-dihydro-1-oxo-1H-inden-5-yl]-2-oxo-5-kernel N-[[3-(2-ethyl-2,CN oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

L23 ANSWER 48 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

1989:423501 HCAPLUS ACCESSION NUMBER:

DOCUMENT NUMBER: 111:23501

Preparation of 5-(aminomethyl)-3-phenyl-2-TITLE:

oxazolidinone derivatives and antibacterial

pharmaceuticals containing them Wang, Chia Lin J.; Wuonola, Mark A. INVENTOR(S): du Pont de Nemours, E. I., and Co., USA PATENT ASSIGNEE(S):

U.S., 13 pp. CODEN: USXXAM SOURCE:

Patent DOCUMENT TYPE: LANGUAGE: English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 4801600	Α	19890131	US 1987-106358	19871009
US 4921869	A	19900501	US 1988-233828	19880819
CA 1322001	A1	19930907	CA 1988-579301	19881004
AU 8823507	A1	19890413	AU 1988-23507	19881006
AU 613669	В2	19910808		
SU 1616518	A3	19901223	SU 1988-4356653	19881006
DK 8805628	Α	19890410	DK 1988-5628	19881007
FI 8804610	Α	19890410	FI 1988-4610	19881007
NO 8804467	Α	19890410	NO 1988-4467	19881007
NO 172890	В	19930614		
NO 172890	С	19930922		
EP 311090	A1	19890412	EP 1988-116621	19881007
R: AT, BE, CH	, DE, E	S, FR, GB,	GR, IT, LI, LU, NL,	
JP 01132569 `	A2	19890525	JP 1988-252207	19881007
ZA 8807550	Α	19900627	ZA 1988-7550	19881007
ни 53359	A2	19901028	HU 1988-5214	19881007
ни 202216	В	19910228		
IL 87972	A1	19930513	IL 1988-87972	19881007
US 4965268	Α	19901023	US 1990-497213	19900315
US 4985429	Α	19910115	US 1990-497211	19900315
US 5032605	Α	19910716	US 1990-497212	19900315
US 5036093	, A	19910730	US 1990-497214	19900315
US 5036092	Α	19910730	US 1990-497215	19900315
US 5039690	Α	19910813	US 1990-497216	19900315
PRIORITY APPLN. INFO.:			US 1987-106358	A3 19871009
			US 1988-233828	A3 19880819

OTHER SOURCE(S):

CASREACT 111:23501; MARPAT 111:23501

GI

$$(CH_2)_n$$
 N
 O
 B
 I

The title compds. I [B = NH2, NR3COR4, NR3S(0)uR5, N3; u = 1, 2; R3 = H,ΑB

alkyl, cycloalkyl; R4 = H, alkyl, alkenyl, cycloalkyl, OR5; R5 = alkyl; X = CH2, O, S, NR6; R6 = H, alkyl; n = 1-3; R1R2 = H2, H and OH, O, H and NR62, NOH, NOR5, NO2CR4, (4-methyl-1-piperazinyl)imino; also when n = 2, then X \neq S; when n = 3, then X \neq O, NR62], specifically the (-)-isomers or the racemic mixts., are prepared for the treatment of bacterial infections. 5-Aminoindan was converted to 5-isocyanatoindan by treatment with HCl gas and phosgene. A solution containing 0.78 g LiBr, 1.96 g Bu3P, and 200 mL xylene was added to a solution containing the latter product

and

18.6 g (R)-glycidyl butyrate to give 26% (-)-[3-(2,3-dihydro-lH-inden-5yl)-2-oxooxazolidin-5-yl]methyl butyrate. The butyrate (39 g) was treated with MeONa to give 82% of the corresponding alc., which (23 g) was mesylated to give 30.8 g mesylate which was subsequently treated with 12.9 g NaN3 to give (-)-5-(azidomethyl)-3-(2,3-dihydro-1H-inden-5-yl)-2oxazolidinone. A mixture containing 400 mL glyme and 14.9 mL (MeO) 3P was added to 25 g of azide; the mixture was heated to 65° for 1 h, and 10 mL 50% HCl were added and the solution was refluxed for 11 h to give the amine-HCl; the latter (8.7 g) was neutralized with NaOH in THF-H2O and treated with 4.14 g Ac2O to give (-)-I (R1R2 = H2, X = CH2, n = 1, B = NHAc) in 100% yield. The latter compound (7 g) was oxidized with CrO3 in aqueous AcOH to give 35% (-)-I (R1R2 = 0, X = CH2, n = 1, B = NHAc) (II). in vitro broth microdilution min. inhibitory concentration of II was 2-4 $\mu g/mL$ and >128 µg/mL against Staphylococcus aureus and Escherichia coli, resp. In an acute lethal mouse model the dose of II required to protect 50% (EC50) of the test animals infected with S. aureus was 1.6 mg/kg. An aqueous suspension for oral administration (5 mL) contains 75 mg active agent, 200 mg Na CM-cellulose, 5 mg NaOBz, 1.0 g sorbitol solution, and 0.025 mL vanillin.

IT 120912-42-7P 120912-43-8P 120912-44-9P 120912-45-0P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of, as bactericide)

RN 120912-42-7 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 120912-43-8 HCAPLUS

CN Acetamide, N-[[3-(2,3-dihydro-1-oxo-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

RN 120912-44-9 HCAPLUS

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CN Acetamide, N-[[3-(2,3-dihydro-1-hydroxy-1H-inden-5-yl)-2-oxo-5-oxazolidinyl]methyl]- (9CI) (CA INDEX NAME)

RN 120912-45-0 HCAPLUS

CN Acetamide, N-[[3-[2,3-dihydro-1-[(4-methyl-1-piperazinyl)imino]-1H-inden-5-yl]-2-oxo-5-oxazolidinyl]methyl]-, (S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

L23 ANSWER 49 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1989:8198 HCAPLUS

DOCUMENT NUMBER:

110:8198

TITLE:

Preparation of (aminomethyl)phenyloxazolidinones as

antibacterial agents

INVENTOR(S):

Gregory, Walter A.

PATENT ASSIGNEE(S):

du Pont de Nemours, E. I., and Co., USA

SOURCE:

U.S., 47 pp. Cont.-in-part of U.S. Ser. No. 676,745,

abandoned.

CODEN: USXXAM

DOCUMENT TYPE:

Patent

LANGUAGE:

English

FAMILY ACC. NUM. COUNT:

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PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE

US 4705799 A 19871110 US 1985-803191 19851202

ZA 8404265 A 19860129 ZA 1984-4265 19840606

HU 196771 IL 77230 CA 1275652 NO 8902178 NO 169122 NO 169122	B A1 A2 A B	19890130 19900610 19901030 19841210 19920203 19920513	HU 1987-5132 IL 1985-77230 CA 1988-580778 NO 1989-2178		19840606 19851204 19881020 19890530
PRIORITY APPLN. INFO.:	C		US 1983-501897 US 1984-578332 US 1984-676745 CA 1984-455844 IL 1984-72028 NO 1984-2273	A2 A2 A3 A	19830607 19840214 19841205 19840605 19840605

OTHER SOURCE(S):

CASREACT 110:8198

GΙ

The title compds. [I; A = NO2, SH, alkylsulfonyl, -sulfinyl, -sulfenyl, etc.; B = N3, (substituted) amino; Y = H, F, Cl, Br, alkyl, NO2; or AY = O(CH2) nO where n = 1, 2, or 3], useful as antibacterial agents for mammals, are prepared A mixture of I (A = 4-MeSO2, B = OSO2C6H4Me-4, Y = H) (preparation given) and NaN3 in DMF was heated at 90-100° for 1 h to give I (A = 4-MeSO2, B = N3, Y = H). = H) (II). II showed a minimal inhibition concentration of 6.3 μ g/mL against Staphylococcus epidermidis.

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of, as antibacterial agent)

RN 96812-87-2 HCAPLUS

CN Acetamide, N-[[3-(1,3-benzodioxol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-(9CI) (CA INDEX NAME)

L23 ANSWER 50 OF 50 HCAPLUS COPYRIGHT 2006 ACS on STN

ACCESSION NUMBER:

1985:437470 HCAPLUS

DOCUMENT NUMBER:

103:37470

TITLE:

Aminomethyloxooxazolidinylbenzene derivatives useful

as antibacterial agents

INVENTOR(S):

Gregory, Walter Adelman

PATENT ASSIGNEE(S):

du Pont de Nemours, E. I., and Co., USA

SOURCE:

Eur. Pat. Appl., 85 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent English

LANGUAGE:

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FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

	KIND DATE	APPLICATION NO.	DATE
EP 127902	A2 19841212		19840605
EP 127902	A3 19870902		
EP 127902	B1 19911016		,
R: AT, BE, CH,	DE, FR, GB, IT,	LI, LU, NL, SE	
ES 533097	A1 19850801	ES 1984-533097	
AU 8429099	A1 19841213	AU 1984-29099	19840605
AU 583250	B2 19890427		
IL 72028	A1 19880531		
CA 1254213	A1 19890516		
AT 68490	E 19911115		
CA 1254213 AT 68490 DK 8402795	A 19841208		
FI 8402273	A 19841208		19840606
FI 83216	B 19910228		
FI 83216	C 19910610		
NO 8402273	A 19841210		19840606
NO 163451	В 19900219		
NO 163451	C 19900530		
JP 60008277			
HU 34462	A2 19850328		19840606
HU 194194	в 19880128		
ZA 8404265			
	В 19890130	ни 1987-5132	. 19840606
SU 1505442			
ES 540812	A1 19880316		
SU 1426451	A3 19880923	SU 1986-4024095	
CA 1275652	A2 19901030		
NO 8902178	A 19841210	NO 1989-2178	19890530
NO 169122	В 19920203		
NO 169122	C 19920513		
PRIORITY APPLN. INFO.:		US 1983-501897	
	•	US 1984-578332	
		CA 1984-455844	
		EP 1984-106400	
		NO 1984-2273	A1 19840606
GI			

$$R^1$$
 N
 CH_2R^2
 I

The bactericidal oxazolidinones I [R = e.g. NO2, cyano, HO, HS, (un)substituted amines, alkylsulfonyl, alkylthio, alkylsulfinyl, aryl, sulfamoyl, alkoxy, or carbamoyl; R1 = H, F, Cl, Br, NO2; RR1 = alkylenedioxy, R2 = NH2, acylamino, N3, alkylsulfonylamino, alkylsulfinylamino] and their physiol. acceptable salts were prepared Thus, (±)-I (R = 4-MeSO2, R1 = H, R2 = Cl) was treated with NaI and the

resulting (\pm) -I (R2 = iodo) treated with NaN3 followed by hydrogenation in F3CCO2H to give (±)-I (R-4-MeSO2; R1 = H, R2 = NH2).F3CCO2H (II). The min. inhibitory concentration of II was 50 μ g/mL against Staphylococcus epidermidis.

96812-87-2P IT

RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of)

96812-87-2 HCAPLUS RN

Acetamide, N-[[3-(1,3-benzodioxol-5-yl)-2-oxo-5-oxazolidinyl]methyl]-CN (CA INDEX NAME)